INTERNATIONAL STANDARD

NORME INTERNATIONALE

ISO 1750

First edition
Première édition
1981-12-15

AMENDMENT 5
AMENDEMENT 5
2008-11-01

Pesticides and other agrochemicals — Common names

AMENDMENT 5

Produits phytosanitaires et assimilés — Noms communs

AMENDEMENT 5

Reference number
Numéro de référence
ISO 1750:1981/Amd.5:2008(E/F)

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Published in Switzerland/Publié en Suisse

Foreword

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International Standards are drafted in accordance with the rules given in the ISO/IEC Directives, Part 2.

The main task of technical committees is to prepare International Standards. Draft International Standards adopted by the technical committees are circulated to the member bodies for voting. Publication as an International Standard requires approval by at least 75 % of the member bodies casting a vote.

Attention is drawn to the possibility that some of the elements of this document may be the subject of patent rights. ISO shall not be held responsible for identifying any or all such patent rights.

Amendment 5 to ISO 1750:1981 was prepared by Technical Committee ISO/TC 81, *Common names for pesticides and other agrochemicals*.

Avant-propos

L'ISO (Organisation internationale de normalisation) est une fédération mondiale d'organismes nationaux de normalisation (comités membres de l'ISO). L'élaboration des Normes internationales est en général confiée aux comités techniques de l'ISO. Chaque comité membre intéressé par une étude a le droit de faire partie du comité technique créé à cet effet. Les organisations internationales, gouvernementales et non gouvernementales, en liaison avec l'ISO participent également aux travaux. L'ISO collabore étroitement avec la Commission électrotechnique internationale (CEI) en ce qui concerne la normalisation électrotechnique.

Les Normes internationales sont rédigées conformément aux règles données dans les Directives ISO/CEI, Partie 2.

La tâche principale des comités techniques est d'élaborer les Normes internationales. Les projets de Normes internationales adoptés par les comités techniques sont soumis aux comités membres pour vote. Leur publication comme Normes internationales requiert l'approbation de 75 % au moins des comités membres votants.

L'attention est appelée sur le fait que certains des éléments du présent document peuvent faire l'objet de droits de propriété intellectuelle ou de droits analogues. L'ISO ne saurait être tenue pour responsable de ne pas avoir identifié de tels droits de propriété et averti de leur existence.

L'Amendement 5 à l'ISO 1750:1981 a été élaboré par le comité technique ISO/TC 81, *Noms communs pour les produits phytosanitaires et assimilés*.

Pesticides and other agrochemicals — Common names

AMENDMENT 5

This fifth Amendment to ISO 1750 supplements the list of common names approved by Technical Committee ISO/TC 81, *Common names for pesticides and other agrochemicals*, for certain pest control chemicals and plant growth regulators of international importance.

In addition to names that have recently been approved, this Amendment includes the names that were approved in Draft Addendum 5 (1984), Draft Addendum 6 (1984), Draft Addendum 7 (1986), and Draft Addendum 8 (1990) but which have not been published.

The common names are listed in alphabetical order in English and are accompanied by the corresponding French name.

The use of each compound is given according to the following classification:

A — Acaricides
AL — Algicides
AT — Attractants
B — Bactericides
F — Fungicides
H — Herbicides
I — Insecticides
IGR — Insect growth regulators
M — Molluscicides
N — Nematicides
P — Plant growth regulators
PA — Plant activators
R — Rodenticides
RE — Repellants
S — Safeners
V — Avicides
Y — Synergists

NOTE 1 When mention is made of more than one use, the letters are arranged alphabetically and not in order of frequency of use.

Produits phytosanitaires et assimilés — Noms communs

AMENDEMENT 5

Le présent cinquième Amendement à l'ISO 1750 complète la liste des noms communs approuvés par le comité technique ISO/TC 81, *Noms communs pour les produits phytosanitaires et assimilés*, pour certains pesticides et autres produits phytopharmaceutiques d'importance internationale.

En complément des noms qui ont été récemment approuvés, le présent Amendement inclut les noms qui ont été approuvés dans le projet d'Additif 5 (1984), le projet d'Additif 6 (1984), le projet d'Additif 7 (1986), et le projet d'Additif 8 (1990) mais qui n'ont pas été publiés.

Les noms communs sont présentés dans l'ordre alphabétique anglais et sont accompagnés du nom commun français correspondant.

L'action de chaque composé est indiquée selon la classification suivante:

A — Acaricides
AL — Algicides
AT — Attractifs
B — Bactéricides
F — Fongicides
H — Herbicides
I — Insecticides
IGR — Régulateurs de la croissance des insectes
M — Molluscicides
N — Nématicides
P — Régulateurs de la croissance des plantes
PA — Activateurs végétaux
R — Rodenticides
RE — Répulsifs
S — Promoteurs de sélectivité
V — Avicides
Y — Synergistes

NOTE 1 Lorsque mention est faite de plus d'une action, les lettres sont disposées par ordre alphabétique et non par ordre de fréquence d'action.

NOTE 2 CAS Registry Number is a Registered Trademark of the American Chemical Society.

NOTE 3 The percentages for mixtures are calculated on a mole fraction basis.

Further amendments to ISO 1750 will be issued in due course giving additional supplementary lists of approved common names. In some cases, widely used names are not available for international use because they are protected by trademarks in some countries.

NOTE 2 Le numéro d'enregistrement CAS est une marque déposée de l'American Chemical Society.

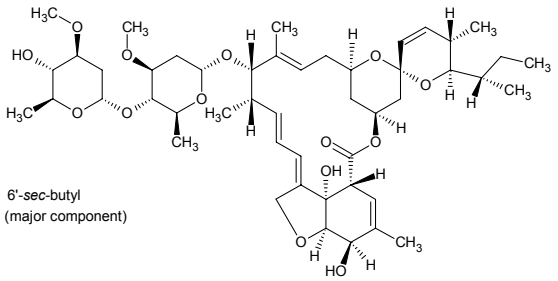
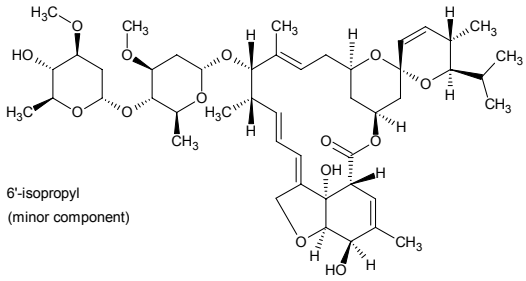
NOTE 3 Les pourcentages des mélanges sont calculés sur une base en fraction molaire.

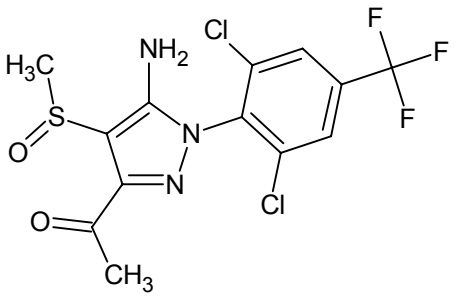
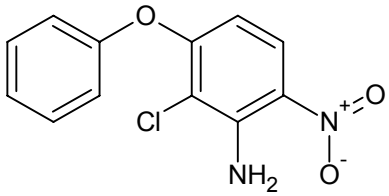
D'autres amendements à l'ISO 1750 sont en cours d'élaboration pour donner des listes supplémentaires de noms communs approuvés. Dans certains cas, des noms largement utilisés ne sont pas acceptables pour un usage international immédiat parce qu'ils sont protégés comme marques commerciales dans certains pays.

1 Approved common names / Noms communs approuvés

1.1 New / Nouveaux

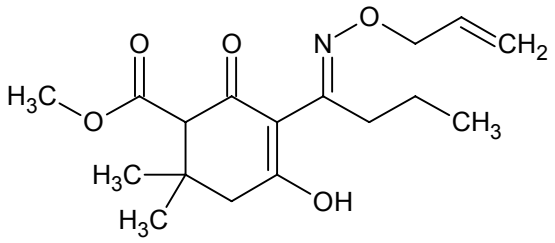
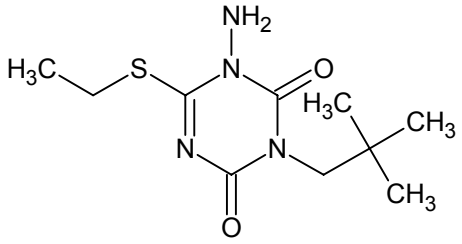
<table>
<tr><th rowspan="2">E: Common name
F: Nom commun</th><th rowspan="2">Chemical name
Nom chimique
E: IUPAC
F: UICPA
C: CAS</th><th colspan="2">Structure
Structure</th><th>Use</th></tr>
<tr><th>Molecular formula
Formule brute</th><th>CAS Registry Number®
Numéro d'enregistrement CAS®</th><th>Applic-ation</th></tr>
<tr><td></td><td colspan="4">IUPAC International Chemical Identifier (InChI™)</td></tr>
<tr><td rowspan="5">E abamectin

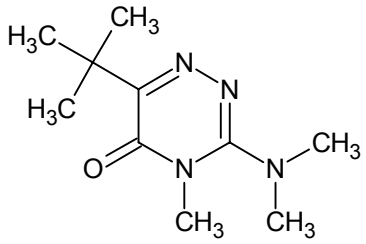
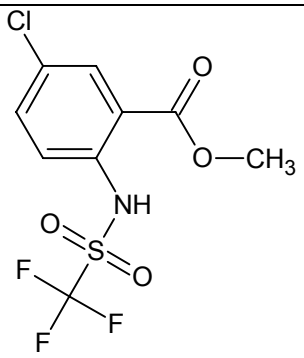
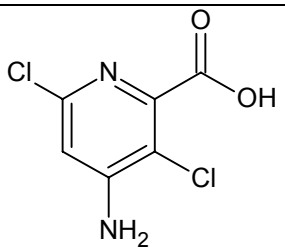
F abamectine
(f)</td><td>mixture of ⩾80 %
(10E,14E,16E)-(1R,4S,5′S,6S,6′R,8R,12S,13S,20R,21R,=24S)-6′-[(S)-sec-butyl]-21,24-dihydroxy-5′,11,13,22-tetramethyl-2-oxo-(3,7,19-trioxatetracyclo=[15.6.1.1^{4,8}.0^{20,24}]pentacosa-10,14,16,22-tetraene)-6-spiro-2′-(5′,6′-dihydro-2′<i>H</i>-pyran)-12-yl 2,6-dideoxy-4-<i>O</i>-(2,6-dideoxy-3-<i>O</i>-methyl-α-L-<i>arabino</i>-hexopyranosyl)-3-<i>O</i>-methyl-α-L-<i>arabino</i>-hexopyranoside
and ⩽20 %
(10<i>E</i>,14<i>E</i>,16<i>E</i>)-(1<i>R</i>,4<i>S</i>,5′<i>S</i>,6<i>S</i>,6′<i>R</i>,8<i>R</i>,12<i>S</i>,13<i>S</i>,20<i>R</i>,21<i>R</i>,=24<i>S</i>)-21,24-dihydroxy-6′-isopropyl-5′,11,13,22-tetramethyl-2-oxo-(3,7,19-trioxatetracyclo[15.6.1.1^{4,8}.0^{20,24}]pentaco=sa-10,14,16,22-tetraene)-6-spiro-2′-(5′,6′-dihydro-2′H-pyran)-12-yl 2,6-dideoxy-4-O-(2,6-dideoxy-3-O-methyl-α-L-arabino-hexopyranosyl)-3-O-methyl-α-L-arabino-hexopyranoside</td><td colspan="2" rowspan="3"> 
6'-sec-butyl
(major component)
 
6'-isopropyl
(minor component)</td><td rowspan="5">A
I
N</td></tr>
<tr><td>mélange de ⩾80 % de
(10E,14E,16E)-(1R,4S,5′S,6S,6′R,8R,12S,13S,20R,21R,=24S)-6′-(S)-sec-butyl-7[[2,6-didésoxy-O-(2,6-didésoxy)-3-O-méthyl-α-L-arabino-hexopyranosyl]oxy]-20,20b-dihydroxy-5′,6,8,19-tétraméthyl-5′,6,6′,7,10,11,14,15,17a,20,20a,20b-dodécahydro-spiro[11.15 méthano-(2H,13H,17H)-furo{4,3.2-pq-[benzodioxacycloocta-2,6-décine]-13 :2′-(2H)-pyran]-17-one (avermectine B$_{1a}$)
et de ⩽20 % de
(10E,14E,16E)-(1R,4S,5′S,6S,6′R,8R,12S,13S,20R,21R,=24S)-6′-isopropyl-7[[2,6-didésoxy-O-(2,6-didésoxy)-3-O-méthyl-α-L-arabino-hexopyranosyl]oxy]-20,20b-dihydroxy-5′,6,8,19-tétraméthyl-5′,6,6′,7,10,11,14,15,17a,20,20a,20b-dodécahydro-spiro[11.15 méthano-(2H,13H,17H)-furo{4,3.2-pq-[benzodioxacycloocta-2,6-décine]-13 :2′-(2H)-pyran]-17-one (avermectine B$_{1b}$)</td></tr>
<tr><td>avermectin B$_1$</td><td>C$_{48}$H$_{72}$O$_{14}$ (avermectin B$_{1a}$)
+
C$_{47}$H$_{70}$O$_{14}$ (avermectin B$_{1b}$)</td><td>71751-41-2</td></tr>
<tr><td colspan="4">avermectin B$_{1a}$ (6′-(S)-sec-butyl)
InChI=1/C48H72O14/c1-11-25(2)43-28(5)17-18-47(62-43)23-34-20-33(61-47)16-15-27(4)42(26(3)13-12-14-32-24-55-45-40(49)29(6)19-35(46(51)58-34)48(32,45)52)59-39-22-37(54-10)44(31(8)57-39)60-38-21-36(53-9)41(50)30(7)56-38/h12-15,17-19,25-26,28,30-31,33-45,49-50,52H,11,16,20-24H2,1-10H3/b13-12+,27-15+,32-14+/t25-,26-,28-,30-,31-,33+,34-,35-,36-,37-,38-,39-,40+,41-,42-,43+,44-,45+,47+,48+/m0/s1</td></tr>
</table>

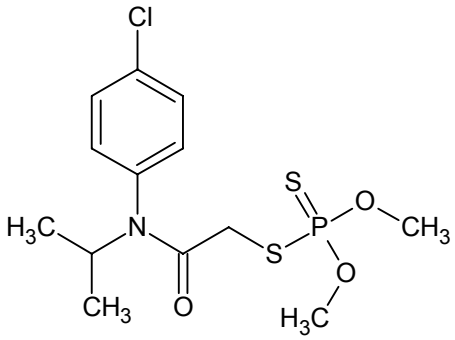
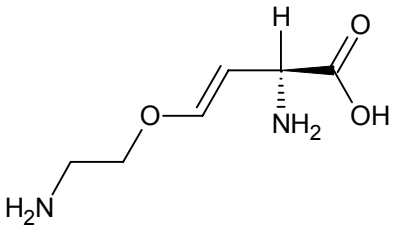
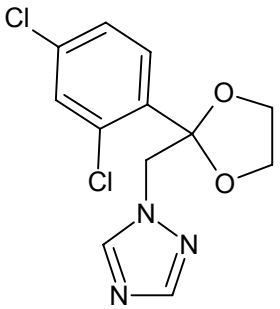
E: Common name F: Nom commun	Chemical name Nom chimique E: IUPAC F: UICPA C: CAS	Structure Structure Molecular formula Formule brute	CAS Registry Number® Numéro d'enregistrement CAS®	Use Applic-ation
	IUPAC International Chemical Identifier (InChI™)			
	avermectin B_{1b} (6′-isopropyl) InChI=1/C47H70O14/c1-24(2)41-27(5)16-17-46(61-41)22-33-19-32(60-46)15-14-26(4)42(25(3)12-11-13-31-23-54-44-39(48)28(6)18-34(45(50)57-33)47(31,44)51)58-38-21-36(53-10)43(30(8)56-38)59-37-20-35(52-9)40(49)29(7)55-37/h11-14,16-18,24-25,27,29-30,32-44,48-49,51H,15,19-23H2,1-10H3/b12-11+,26-14+,31-13+/t25-,27-,29-,30-,32+,33-,34-,35-,36-,37-,38-,39+,40-,41+,42-,43-,44+,46+,47+/m0/s1		
E acetoprole **F** acétoprole (m)	1-[5-amino-1-(2,6-dichloro-α,α,α-trifluoro-*p*-tolyl)-4-(methylsulfinyl)pyrazol-3-yl]ethanone 1-[5-amino-1-(2,6-dichloro-4-trifluorométhylphényl)-4-(méthylsulfinyl)pyrazol-3-yl]éthanone 1-[5-amino-1-[2,6-dichloro-4-(trifluoromethyl)phenyl]-4-(methylsulfinyl)-1*H*-pyrazol-3-yl]= ethanone			**A I N**
		$C_{13}H_{10}Cl_2F_3N_3O_2S$	209861-58-5	
	InChI=1/C13H10Cl2F3N3O2S/c1-5(22)9-11(24(2)23)12(19)21(20-9)10-7(14)3-6(4-8(10)15)13(16,17)18/h3-4H,19H2,1-2H3			
E aclonifen **F** aclonifène (m)	2-chloro-6-nitro-3-phenoxyaniline 2-chloro-6-nitro-3-phénoxyaniline 2-chloro-6-nitro-3-phenoxybenzenamine			**H**
		$C_{12}H_9ClN_2O_3$	74070-46-5	
	InChI=1/C12H9ClN2O3/c13-11-10(18-8-4-2-1-3-5-8)7-6-9(12(11)14)15(16)17/h1-7H,14H2			
	NOTE The name “aclonifen” is not acceptable for use in France because of the risk of confusion with the WHO name “clomifen”. NOTE Le nom «aclonifène» n'est pas acceptable pour l'emploi en France, car il entre en conflit avec le nom OMS «clomifène».			

<table>
<tr><th rowspan="3">E: Common name
F: Nom commun</th><th rowspan="3">Chemical name
Nom chimique
E: IUPAC
F: UICPA
C: CAS</th><th colspan="2">Structure
Structure</th><th>Use</th></tr>
<tr><th>Molecular formula
Formule brute</th><th>CAS Registry Number®
Numéro d'enregistrement CAS®</th><th>Applic-ation</th></tr>
<tr><td colspan="4">IUPAC International Chemical Identifier (InChI™)</td></tr>
<tr><td rowspan="5">E acypetacs

F acypétacs (m)</td><td>mixture of C_8 to C_{10} linear and branched chain saturated aliphatic carboxylic acids, the branched chain acids predominating and consisting of:
a) acids in which the main chain is dialkyl-substituted on the second carbon atom; and
b) acids in which the second carbon atom is either unsubstituted or monoalkyl-substituted.
Both (a) and (b) acids may be further alkyl-substituted on the third or higher carbon atoms</td><td colspan="2" rowspan="3">n/a</td><td rowspan="5">F</td></tr>
<tr><td>mélange d'acides carboxyliques aliphatiques saturés, linéaires ou ramifiés, en C_8 à C_{10}, dans lequel les acides à chaîne ramifiée prédominent et sont des types suivants:
a) acides dont la chaîne principale comporte deux substituants alkyles sur le deuxième atome de carbone;
b) acides dans lesquels le deuxième atome de carbone est soit non substitué soit substitué par un seul alkyle.
Les atomes de carbone en position 3 ou supérieure des acides a) et b) peuvent être substitués par d'autres alkyles.</td></tr>
<tr><td>acypetacs</td></tr>
<tr><td>n/a</td><td></td></tr>
<tr><td colspan="3">NOTE It should be stated which salt is present, for example acypetacs-copper or acypetacs-zinc [380221-54-5].
NOTE Il convient de préciser quel est le sel présent, par exemple acypétacs-cuivre ou acypétacs-zinc [380221-54-5].</td></tr>
</table>

E: Common name F: Nom commun	Chemical name Nom chimique E: IUPAC F: UICPA C: CAS	Structure Structure Molecular formula Formule brute	CAS Registry Number® Numéro d'enregistrement CAS®	Use Application
	IUPAC International Chemical Identifier (InChI™)			
E alloxydim **F** alloxydim (m)	methyl (1*RS*)-3-[(*E*)-1-(allyloxyimino)butyl]-4-hydroxy-6,6-dimethyl-2-oxocyclohex-3-enecarboxylate (1*RS*)-3-[(*E*)-1-(allyloxy)imino]butyl]-4-hydroxy-6,6-diméthyl-2-oxocyclohex-3-ène-1-carboxylate de méthyle methyl 2,2-dimethyl-4,6-dioxo-5-[(1*E*)-1-[(2-propenyloxy)imino]=butyl]cyclohexanecarboxylate			**H**
		$C_{17}H_{25}NO_5$	55634-91-8	
	InChI=1/C17H25NO5/c1-6-8-11(18-23-9-7-2)13-12(19)10-17(3,4)14(15(13)20)16(21)22-5/h7,14,19H,2,6,8-10H2,1,3-5H3/b18-11+/t14-/s3 CAS-preferred tautomer InChI=1/C17H25NO5/c1-6-8-11(18-23-9-7-2)13-12(19)10-17(3,4)14(15(13)20)16(21)22-5/h7,13-14H,2,6,8-10H2,1,3-5H3/b18-11+/t13?,14? alloxydim-sodium InChI=1/C17H25NO5.Na/c1-6-8-11(18-23-9-7-2)13-12(19)10-17(3,4)14(15(13)20)16(21)22-5;/h7,14,19H,2,6,8-10H2,1,3-5H3;/q;+1/p-1/b18-11+;/t14-;/s3/fC17H24NO5.Na/h19h;/q-1;m			
	NOTE 1 It should be stated which salt is present, for example alloxydim-sodium [55635-13-7]. NOTE 1 Il convient de préciser quel est le sel présent, par exemple alloxydim-sodium [55635-13-7]. NOTE 2 The CAS name and Registry Number are for the tautomer that is preferred under the CAS nomenclature rules. NOTE 2 Le nom et le numéro de registre CAS sont attribués au tautomère préféré conformément aux règles de nomenclature CAS.			
E ametridione **F** amétridione (f)	1-amino-6-ethylthio-3-neopentyl-1,3,5-triazine-2,4(1*H*,3*H*)-dione 1-amino-6-éthylthio-3-néopentyl-1,3,5-triazine-(1*H*,3*H)*-2,4-dione 1-amino-3-(2,2-dimethylpropyl)-6-(ethylthio)-1,3,5-triazine-2,4(1*H*,3*H*)-dione			**H**
		$C_{10}H_{18}N_4O_2S$	78168-93-1	
	InChI=1/C10H18N4O2S/c1-5-17-8-12-7(15)13(6-10(2,3)4)9(16)14(8)11/h5-6,11H2,1-4H3			

E: Common name F: Nom commun	Chemical name Nom chimique E: IUPAC F: UICPA C: CAS	Structure Structure Molecular formula Formule brute	CAS Registry Number® Numéro d'enregistrement CAS®	Use Application
	IUPAC International Chemical Identifier (InChI™)			
E amibuzin **F** amibuzine (f)	6-*tert*-butyl-3-dimethylamino-4-methyl-1,2,4-triazin-5(4*H*)-one 6-*tert*-butyl-3-diméthylamino-4-méthyl-(4*H*)-1,2,4-triazin-5-one 3-(dimethylamino)-6-(1,1-dimethylethyl)-4-methyl-1,2,4-triazin-5(4*H*)-one			**H**
		$C_{10}H_{18}N_4O$	76636-10-7	
	InChI=1/C10H18N4O/c1-10(2,3)7-8(15)14(6)9(12-11-7)13(4)5/h1-6H3			
	NOTE 1 The name “amibuzin” is not acceptable for use in France because of the risk of confusion with the trade name “Aminozine”. NOTE 1 Le nom «amibuzine» n'est pas acceptable pour l'emploi en France, car il entre en conflit avec le nom commercial «Aminozine». NOTE 2 The name “amibuzin” is not acceptable for use in Japan because of the risk of confusion with the trade name “Triamidine”. NOTE 2 Le nom «amibuzine» n'est pas acceptable pour l'emploi au Japon, car il entre en conflit avec le nom commercial «Triamidine».			
E amidoflumet **F** amidoflumet (m)	methyl 5-chloro-2-{[(trifluoromethyl)=sulfonyl]amino}benzoate 5-chloro-2-{[(trifluorométhyl)=sulfonyl]amino}benzoate de méthyle methyl 5-chloro-2-[[(trifluoromethyl)=sulfonyl]amino]benzoate			**A**
		$C_9H_7ClF_3NO_4S$	84466-05-7	
	InChI=1/C9H7ClF3NO4S/c1-18-8(15)6-4-5(10)2-3-7(6)14-19(16,17)9(11,12)13/h2-4,14H,1H3			
E aminopyralid **F** aminopyralide (m)	4-amino-3,6-dichloropyridine-2-carboxylic acid acide 4-amino-3,6-dichloropyridine-2-carboxylique 4-amino-3,6-dichloro-2-pyridinecarboxylic acid			**H**
		$C_6H_4Cl_2N_2O_2$	150114-71-9	
	InChI=1/C6H4Cl2N2O2/c7-3-1-2(9)4(8)5(10-3)6(11)12/h1H,(H2,9,10)(H,11,12)/f/h11H,9H2 aminopyralid-potassium InChI=1/C6H4Cl2N2O2.K/c7-3-1-2(9)4(8)5(10-3)6(11)12;/h1H,(H2,9,10)(H,11,12);/q;+1/p-1/fC6H3Cl2N2O2.K/h9H2;/q-1;m			
	NOTE It should be stated which ester or salt is present, for example aminopyralid-potassium [566191-87-5]. NOTE Il convient de préciser quel est l'ester ou le sel présent, par exemple aminopyralide-potassium [566191-87-5].			

E: Common name F: Nom commun	Chemical name Nom chimique E: IUPAC F: UICPA C: CAS	Structure Structure Molecular formula Formule brute	CAS Registry Number® Numéro d'enregistrement CAS®	Use Applic-ation
	IUPAC International Chemical Identifier (InChI™)			
E anilofos **F** anilofos (m)	*S*-4-chloro-*N*-isopropylcarbaniloylmethyl *O,O*-dimethyl phosphorodithioate dithiophosphate de *S*-[[*N*-(4-chlorophényl)-*N*-isopropyl-carbamoyl]méthyle] et de *O,O*-diméthyle *S*-[2-[(4-chlorophenyl)(1-methylethyl)amino]-2-oxoethyl] *O,O*-dimethyl phosphorodithioate			**H**
		$C_{13}H_{19}ClNO_3PS_2$	64249-01-0	
	InChI=1/C13H19ClNO3PS2/c1-10(2)15(12-7-5-11(14)6-8-12)13(16)9-21-19(20,17-3)18-4/h5-8,10H,9H2,1-4H3			
	NOTE The name “anilofos” is not acceptable for use in Japan because of the risk of confusion with the trade name “Amilofos”. NOTE Le nom «anilofos» n'est pas acceptable pour l'emploi au Japon, car il entre en conflit avec le nom commercial «Amilofos».			
E aviglycine **F** aviglycine (f)	(*E*)-L-2-[2-(2-aminoethoxy)vinyl]glycine (*E*)-L-2-[2-(2-aminoéthoxy)vinyl]glycine (2*S*,3*E*)-2-amino-4-(2-aminoethoxy)-3-butenoic acid			**P**
		$C_6H_{12}N_2O_3$	49669-74-1	
	InChI=1/C6H12N2O3/c7-2-4-11-3-1-5(8)6(9)10/h1,3,5H,2,4,7-8H2,(H,9,10)/b3-1+/t5-/m0/s1/f/h9H aviglycine hydrochloride InChI=1/C6H12N2O3.ClH/c7-2-4-11-3-1-5(8)6(9)10;/h1,3,5H,2,4,7-8H2,(H,9,10);1H/b3-1+;/t5-;/m0./s1/f/h9H;			
	NOTE It should be stated which salt is present, for example aviglycine hydrochloride [55720-26-8]. NOTE Il convient de préciser quel est le sel présent, par exemple le chlorhydrate d'aviglycine [55720-26-8].			
E azaconazole **F** azaconazole (m)	1-{[2-(2,4-dichlorophenyl)-1,3-dioxolan-2-yl]methyl}-1*H*-1,2,4-triazole 1-{[2-(2,4-dichlorophényl)-1,3-dioxolan-2-yl]méthyl}-1*H*-1,2,4-triazole 1-[[2-(2,4-dichlorophenyl)-1,3-dioxolan-2-yl]methyl]-1*H*-1,2,4-triazole			**F**
		$C_{12}H_{11}Cl_2N_3O_2$	60207-31-0	
	InChI=1/C12H11Cl2N3O2/c13-9-1-2-10(11(14)5-9)12(18-3-4-19-12)6-17-8-15-7-16-17/h1-2,5,7-8H,3-4,6H2			

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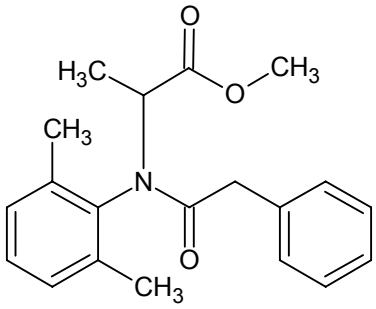
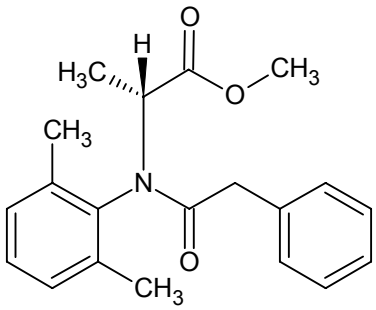
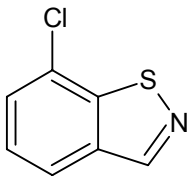
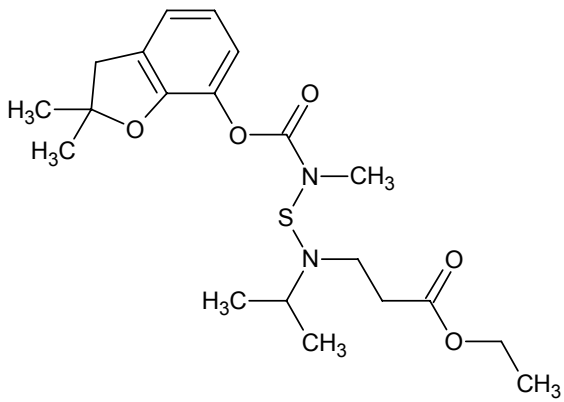
	IUPAC International Chemical Identifier (InChI™)			
E benalaxyl **F** bénalaxyl (m)	methyl *N*-(phenylacetyl)-*N*-(2,6-xylyl)-DL-alaninate DL-(*N*-phénylacétyl-*N*-2,6-xylyl)alaninate de méthyle methyl *N*-(2,6-dimethylphenyl)-*N*-(phenylacetyl)-DL-alaninate			**F**
		$C_{20}H_{23}NO_3$	71626-11-4	
	InChI=1/C20H23NO3/c1-14-9-8-10-15(2)19(14)21(16(3)20(23)24-4)18(22)13-17-11-6-5-7-12-17/h5-12,16H,13H2,1-4H3/t16-/s3			
E benalaxyl-M **F** bénalaxyl-M (m)	methyl *N*-(phenylacetyl)-*N*-(2,6-xylyl)-D-alaninate D-(*N*-phénylacétyl-*N*-2,6-xylyl)alaninate de méthyle methyl *N*-(2,6-dimethylphenyl)-*N*-(phenylacetyl)-D-alaninate			**F**
		$C_{20}H_{23}NO_3$	98243-83-5	
	InChI=1/C20H23NO3/c1-14-9-8-10-15(2)19(14)21(16(3)20(23)24-4)18(22)13-17-11-6-5-7-12-17/h5-12,16H,13H2,1-4H3/t16-/m1/s1			
E benclothiaz **F** benclothiaze (m)	7-chloro-1,2-benzothiazole 7-chloro-1,2-benzisothiazole 7-chloro-1,2-benzisothiazole			**N**
		C_7H_4ClNS	89583-90-4	
	InChI=1/C7H4ClNS/c8-6-3-1-2-5-4-9-10-7(5)6/h1-4H			
E benfuracarb **F** benfuracarb (m)	ethyl *N*-[2,3-dihydro-2,2-dimethylbenzofuran-7-yloxycarbonyl(methyl)aminothio]-*N*-isopropyl-β-alaninate *N*-[2,3-dihydro-2,2-diméthylbenzofuran-7-yloxycarbonyl(méthylamino)thio]-*N*-isopropyl-β-alaninate d'éthyle 2,3-dihydro-2,2-dimethyl-7-benzofuranyl 2-methyl-4-(1-methylethyl)-7-oxo-8-oxa-3-thia-2,4-diazadecanoate			**I**
		$C_{20}H_{30}N_2O_5S$	82560-54-1	
	InChI=1/C20H30N2O5S/c1-7-25-17(23)11-12-22(14(2)3)28-21(6)19(24)26-16-10-8-9-15-13-20(4,5)27-18(15)16/h8-10,14H,7,11-13H2,1-6H3			

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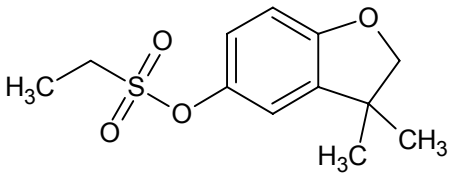
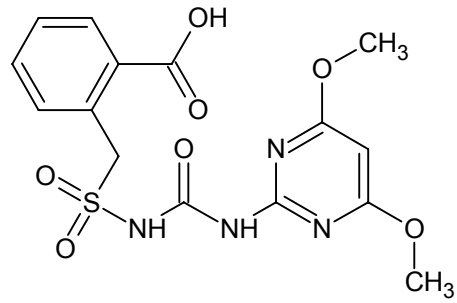
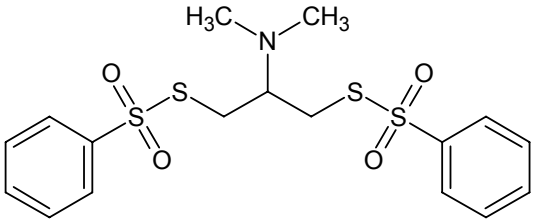
E benfuresate **F** benfurésate (m)	2,3-dihydro-3,3-dimethylbenzofuran-5-yl ethanesulfonate éthanesulfonate de 2,3-dihydro-3,3-diméthylbenzofuran-5-yle 2,3-dihydro-3,3-dimethyl-5-benzofuranyl ethanesulfonate			**H**
		$C_{12}H_{16}O_4S$	68505-69-1	
	InChI=1/C12H16O4S/c1-4-17(13,14)16-9-5-6-11-10(7-9)12(2,3)8-15-11/h5-7H,4,8H2,1-3H3			
E bensulfuron **F** bensulfuron (m)	α-[(4,6-dimethoxypyrimidin-2-ylcarbamoyl)sulfamoyl]-*o*-toluic acid acide 2-[(4,6-diméthoxypyrimidin-2-yl)-3-uréidosulfonylméthyl]benzoïque 2-[[[[[(4,6-dimethoxy-2-pyrimidinyl)amino]carbonyl]amino]= sulfonyl]methyl]benzoic acid			**H**
		$C_{15}H_{16}N_4O_7S$	99283-01-9	
	InChI=1/C15H16N4O7S/c1-25-11-7-12(26-2)17-14(16-11)18-15(22)19-27(23,24)8-9-5-3-4-6-10(9)13(20)21/h3-7H,8H2,1-2H3,(H,20,21)(H2,16,17,18,19,22)/f/h18-20H bensulfuron-methyl InChI=1/C16H18N4O7S/c1-25-12-8-13(26-2)18-15(17-12)19-16(22)20-28(23,24)9-10-6-4-5-7-11(10)14(21)27-3/h4-8H,9H2,1-3H3,(H2,17,18,19,20,22)/f/h19-20H			
	NOTE It should be stated which ester or salt is present, for example bensulfuron-methyl [83055-99-6]. NOTE Il convient de préciser quel est l'ester ou le sel présent, par exemple bensulfuron-méthyle [83055-99-6].			
E bensultap **F** bensultap (m)	*S,S*′-2-dimethylaminotrimethylene di(benzenethiosulfonate) dibenzènethiosulfonate de *S,S*′-[2-(diméthylamino)]propan-1,3-diyle *S,S*′-[2-(dimethylamino)-1,3-propanediyl] di(benzenesulfonothioate)			**I**
		$C_{17}H_{21}NO_4S_4$	17606-31-4	
	InChI=1/C17H21NO4S4/c1-18(2)15(13-23-25(19,20)16-9-5-3-6-10-16)14-24-26(21,22)17-11-7-4-8-12-17/h3-12,15H,13-14H2,1-2H3			

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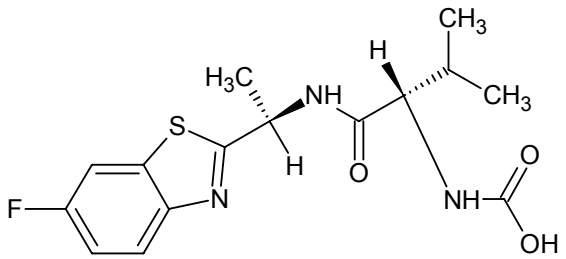
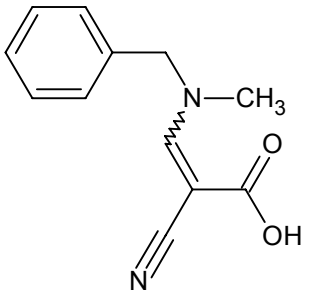
E benthiavalicarb **F** benthiavalicarbe (m)	[(*S*)-1-{[(1*R*)-1-(6-fluoro-1,3-benzothiazol-2-yl)ethyl]carbamoyl}-2-methylpropyl]carbamic acid acide [(*S*)-1-{[(1*R*)-1-(6-fluoro-1,3-benzothiazol-2-yl)éthyl]carbamoyl}-2-méthylpropyl]carbamique [(1*S*)-1-[[[(1*R*)-1-(6-fluoro-2-benzothiazolyl)ethyl]amino]carbonyl]-2-methylpropyl]carbamic acid			**F**
		$C_{15}H_{18}FN_3O_3S$	413615-35-7	
	InChI=1/C15H18FN3O3S/c1-7(2)12(19-15(21)22)13(20)17-8(3)14-18-10-5-4-9(16)6-11(10)23-14/h4-8,12,19H,1-3H3,(H,17,20)(H,21,22)/t8-,12+/m1/s1/f/h17,21H benthiavalicarb-isopropyl InChI=1/C18H24FN3O3S/c1-9(2)15(22-18(24)25-10(3)4)16(23)20-11(5)17-21-13-7-6-12(19)8-14(13)26-17/h6-11,15H,1-5H3,(H,20,23)(H,22,24)/t11-,15+/m1/s1/f/h20,22H			
	NOTE It should be stated which ester or salt is present, for example benthiavalicarb-isopropyl [177406-68-7]. NOTE Il convient de préciser quel est l'ester ou le sel présent, par exemple le benthiavalicarbe-isopropyle [177406-68-7].			
E benzamacril **F** benzamacril (m)	(*EZ*)-3-[benzyl(methyl)amino]-2-cyanoacrylic acid acide (*EZ*)-3-[méthyl(benzyl)amino]-2-cyanoacrylique 2-cyano-3-[methyl(phenylmethyl)amino]-2-propenoic acid			**F**
		$C_{12}H_{12}N_2O_2$	127087-86-9	
	InChI=1/C12H12N2O2/c1-14(9-11(7-13)12(15)16)8-10-5-3-2-4-6-10/h2-6,9H,8H2,1H3,(H,15,16)/b11-9?/f/h15H benzamacril-isobutyl InChI=1/C16H20N2O2/c1-13(2)12-20-16(19)15(9-17)11-18(3)10-14-7-5-4-6-8-14/h4-8,11,13H,10,12H2,1-3H3/b15-11?			
	NOTE 1 It should be stated which ester or salt is present, for example benzamacril-isobutyl [88107-27-1]. NOTE 1 Il convient de préciser quel est l'ester ou le sel présent, par exemple benzamacril-isobutyle [88107-27-1]. NOTE 2 The (*E*)- and (*Z*)-isomer population is temperature dependent. NOTE 2 La proportion des isomères (*E*) et (*Z*) dépend de la température.			

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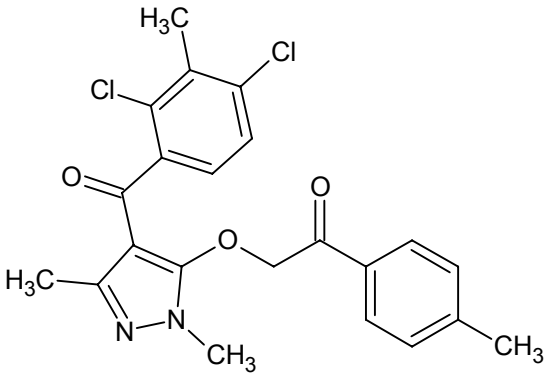
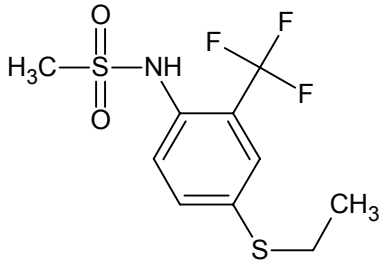
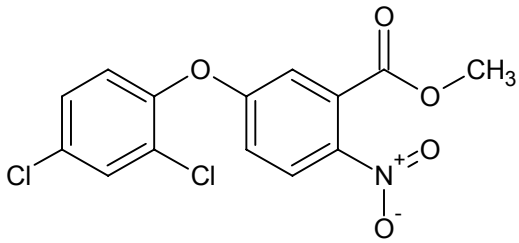
E benzofenap **F** benzofénap (m)	2-[4-(2,4-dichloro-*m*-toluoyl)-1,3-dimethylpyrazol-5-yloxy]-4′-methylacetophenone 2-[4-(2,4-dichloro-3-méthylbenzoyl)-1,3-diméthyl-(1*H*)-pyrazol-5-yloxy-*p*-tolylacétone 2-[[4-(2,4-dichloro-3-methylbenzoyl)-1,3-dimethyl-1*H*-pyrazol-5-yl]oxy]-1-(4-methylphenyl)ethanone			**H**
		$C_{22}H_{20}Cl_2N_2O_3$	82692-44-2	
	InChI=1/C22H20Cl2N2O3/c1-12-5-7-15(8-6-12)18(27)11-29-22-19(14(3)25-26(22)4)21(28)16-9-10-17(23)13(2)20(16)24/h5-10H,11H2,1-4H3			
E benzofluor **F** benzofluor (m)	4′-ethylthio-2′-(trifluoromethyl)methylsulfonanilide *N*-[4-éthylthio-2-(trifluorométhyl)-phényl]méthanesulfonamide *N*-[4-(ethylthio)-2-(trifluoromethyl)-phenyl]methanesulfonamide			**H** **P**
		$C_{10}H_{12}F_3NO_2S_2$	68672-17-3	
	InChI=1/C10H12F3NO2S2/c1-3-17-7-4-5-9(14-18(2,15)16)8(6-7)10(11,12)13/h4-6,14H,3H2,1-2H3			
E bifenox **F** bifénox (m)	methyl 5-(2,4-dichlorophenoxy)-2-nitrobenzoate 5-(2,4-dichlorophénoxy)-2-nitrobenzoate de méthyle methyl 5-(2,4-dichlorophenoxy)-2-nitrobenzoate			**H**
		$C_{14}H_9Cl_2NO_5$	42576-02-3	
	InChI=1/C14H9Cl2NO5/c1-21-14(18)10-7-9(3-4-12(10)17(19)20)22-13-5-2-8(15)6-11(13)16/h2-7H,1H3			

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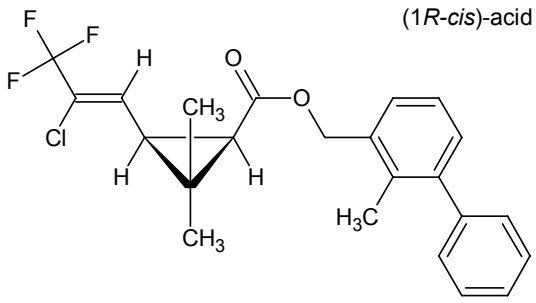
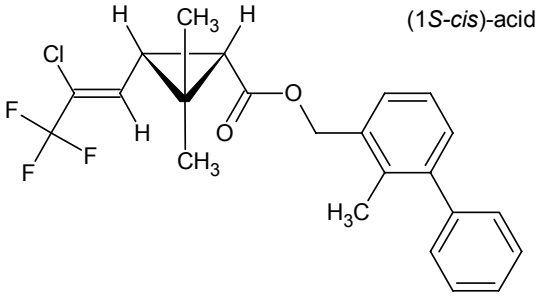
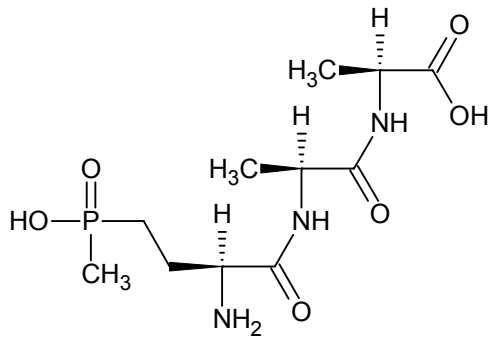
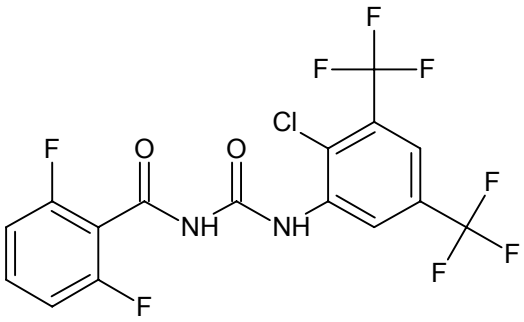
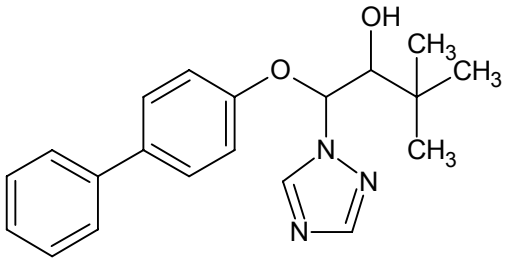
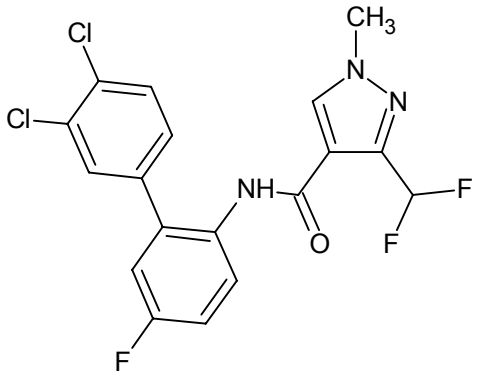
	IUPAC International Chemical Identifier (InChI™)			
E bifenthrin **F** bifenthrine (f)	2-methylbiphenyl-3-ylmethyl (1*RS*,3*RS*)-3-[(*Z*)-2-chloro-3,3,3-trifluoroprop-1-enyl]-2,2-dimethylcyclopropanecarboxylate Alternative Rothamsted-style stereodescriptors: 2-methylbiphenyl-3-ylmethyl (1*RS*)-*cis*-3-[(*Z*)-2-chloro-3,3,3-trifluoroprop-1-enyl]-2,2-dimethylcyclopropanecarboxylate			**A** **I**
	(1*RS*,3*RS*)-3-[(*Z*)-2-chloro-3,3,3-trifluoroprop-1-ènyl]-2,2-diméthylcyclopropanecarboxylate de 2-méthylbiphényl-3-ylméthyle			
	(2-methyl[1,1′-biphenyl]-3-yl)methyl (1*R*,3*R*)-*rel*-3-[(1*Z*)-2-chloro-3,3,3-trifluoro-1-propenyl]-2,2-dimethylcyclopropanecarboxylate	$C_{23}H_{22}ClF_3O_2$	82657-04-3	
	InChI=1/C23H22ClF3O2/c1-14-11-15(9-10-17(14)16-7-5-4-6-8-16)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-/s3			
E bilanafos **F** bilanafos (m)	(2*S*)-2-amino-4-[hydroxy(methyl)=phosphinoyl]butyryl-L-alanyl-L-alanine *N*-(*N*-((2*S*)-2-amino-4-(hydroxy(méthyl)phosphinoyl)butyryl)-L-alanyl)-L-alanine 4-(hydroxymethylphosphinyl)-L-2-aminobutanoyl-L-alanyl-L-alanine			**H**
		$C_{11}H_{22}N_3O_6P$	35597-43-4	
	InChI=1/C11H22N3O6P/c1-6(9(15)14-7(2)11(17)18)13-10(16)8(12)4-5-21(3,19)20/h6-8H,4-5,12H2,1-3H3,(H,13,16)(H,14,15)(H,17,18)(H,19,20)/t6-,7-,8-/m0/s1/f/h13-14,17,19H bilanafos-sodium InChI=1/C11H22N3O6P.Na/c1-6(9(15)14-7(2)11(17)18)13-10(16)8(12)4-5-21(3,19)20;/h6-8H,4-5,12H2,1-3H3,(H,13,16)(H,14,15)(H,17,18)(H,19,20);/q;+1/p-1/t6-,7-,8-;/m0./s1/fC11H21N3O6P.Na/h13-14,17H;/q-1;m			
	NOTE It should be stated which ester or salt is present, for example bilanafos-sodium [71048-99-2]. NOTE Il convient de préciser quel est l'ester ou le sel présent, par exemple bilanafos-sodium [71048-99-2].			

E: Common name F: Nom commun	Chemical name Nom chimique E: IUPAC F: UICPA C: CAS	Structure Structure Molecular formula Formule brute	CAS Registry Number® Numéro d'enregistrement CAS®	Use Applic-ation
	IUPAC International Chemical Identifier (InChI™)			
E bistrifluron **F** bistrifluron (m)	1-[2-chloro-3,5-bis(trifluoromethyl)phenyl]-3-(2,6-difluorobenzoyl)urea 1-[2-chloro-3,5-bis(trifluorométhyl)phényl]-3-(2,6-difluorobenzoyl)urée *N*-[[[2-chloro-3,5-bis(trifluoromethyl)=phenyl]amino]carbonyl]-2,6-difluorobenzamide			**IGR**
		$C_{16}H_7ClF_8N_2O_2$	201593-84-2	
	InChI=1/C16H7ClF8N2O2/c17-12-7(16(23,24)25)4-6(15(20,21)22)5-10(12)26-14(29)27-13(28)11-8(18)2-1-3-9(11)19/h1-5H,(H2,26,27,28,29)/f/h26-27H			
E bitertanol **F** bitertanol (m)	(1*RS*,2*RS*;1*RS*,2*SR*)-1-(biphenyl-4-yloxy)-3,3-dimethyl-1-(1*H*-1,2,4-triazol-1-yl)butan-2-ol (20:80 ratio of (1*RS*,2*RS*)- and (1*RS*,2*SR*)-isomers) 1-(biphényl-4-yloxy]-1-((1*H*)1,2,4-triazol-1-yl)-3,3-diméthylbutan-2-ol β-([1,1′-biphenyl]-4-yloxy)-α-(1,1-dimethylethyl)-1*H*-1,2,4-triazole-1-ethanol			**F**
		$C_{20}H_{23}N_3O_2$	55179-31-2	
	InChI=1/C20H23N3O2/c1-20(2,3)18(24)19(23-14-21-13-22-23)25-17-11-9-16(10-12-17)15-7-5-4-6-8-15/h4-14,18-19,24H,1-3H3			
E bixafen **F** bixafène (m)	*N*-(3′,4′-dichloro-5-fluoro[1,1′-biphenyl]-2-yl)-3-(difluoromethyl)-1-methyl-1*H*-pyrazole-4-carboxamide *N*-(3′,4′-dichloro-5-fluorobiphényl-2-yl)-3-(difluorométhyl)-1-méthyl-1*H*-pyrazole-4-carboxamide *N*-(3′,4′-dichloro-5-fluorobiphenyl-2-yl)-3-(difluoromethyl)-1-methyl-1*H*-pyrazole-4-carboxamide			**F**
		$C_{18}H_{12}Cl_2F_3N_3O$	581809-46-3	
	InChI=1/C18H12Cl2F3N3O/c1-26-8-12(16(25-26)17(22)23)18(27)24-15-5-3-10(21)7-11(15)9-2-4-13(19)14(20)6-9/h2-8,17H,1H3,(H,24,27)/f/h24H			

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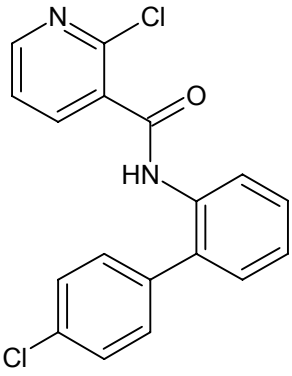
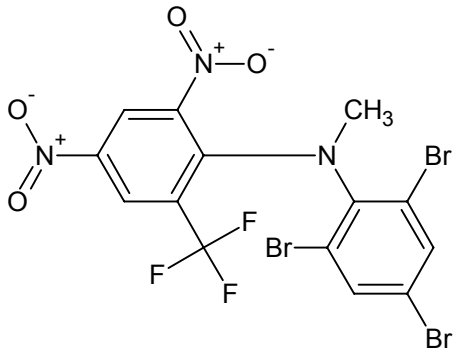
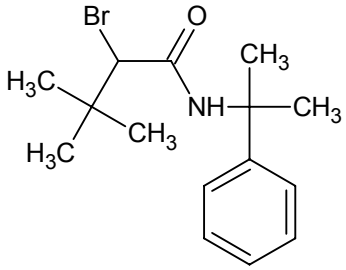
	IUPAC International Chemical Identifier (InChI™)			
E boscalid **F** boscalide (m)	2-chloro-*N*-(4′-chlorobiphenyl-2-yl)nicotinamide 2-chloro-*N*-(4′-chlorobiphényl-2-yl)nicotinamide 2-chloro-*N*-(4′-chloro[1,1′-biphenyl]-2-yl)-3-pyridinecarboxamide			**F**
		$C_{18}H_{12}Cl_2N_2O$	188425-85-6	
	InChI=1/C18H12Cl2N2O/c19-13-9-7-12(8-10-13)14-4-1-2-6-16(14)22-18(23)15-5-3-11-21-17(15)20/h1-11H,(H,22,23)/f/h22H			
E bromethalin **F** bromét haline (f)	α,α,α-trifluoro-*N*-methyl-4,6-dinitro-*N*-(2,4,6-tribromophenyl)-*o*-toluidine [(2,4-dinitro-6-trifluorométhyl)phényl]méthyl(2,4,6-tribromophényl)amine *N*-methyl-2,4-dinitro-*N*-(2,4,6-tribromophenyl)-6-(trifluoromethyl)benzenamine			**R**
		$C_{14}H_7Br_3F_3N_3O_4$	63333-35-7	
	InChI=1/C14H7Br3F3N3O4/c1-21(13-9(16)2-6(15)3-10(13)17)12-8(14(18,19)20)4-7(22(24)25)5-11(12)23(26)27/h2-5H,1H3			
E bromobutide **F** bromobutide (m)	(*RS*)-2-bromo-3,3-dimethyl-*N*-(1-methyl-1-phenylethyl)butyramide *N*-[(*RS*)-2-phénylprop-2-yl]-2-bromo-3,3-diméthylbutyramide 2-bromo-3,3-dimethyl-*N*-(1-methyl-1-phenylethyl)butanamide			**H**
		$C_{15}H_{22}BrNO$	74712-19-9	
	InChI=1/C15H22BrNO/c1-14(2,3)12(16)13(18)17-15(4,5)11-9-7-6-8-10-11/h6-10,12H,1-5H3,(H,17,18)/t12-/s3/f/h17H			

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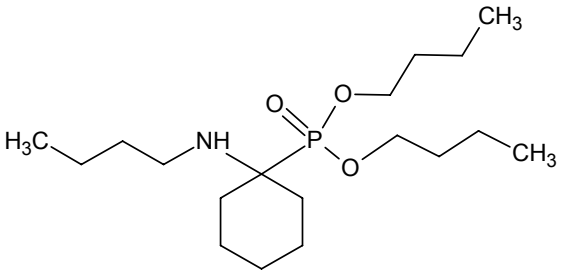
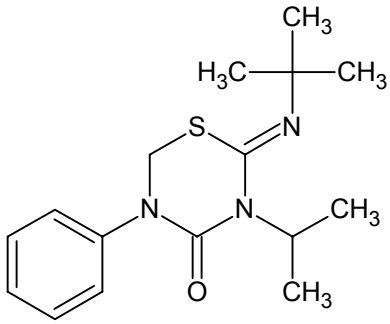
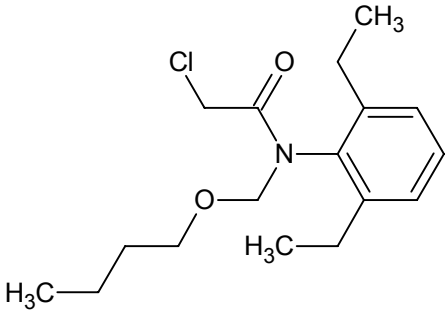
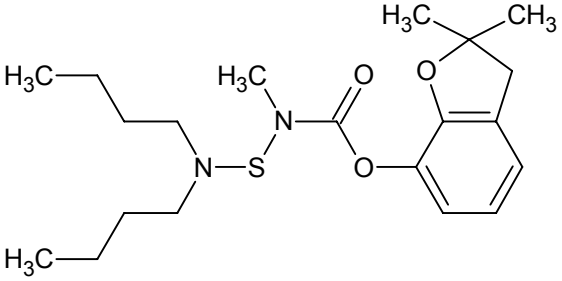
E buminafos **F** buminafos (m)	dibutyl 1-butylaminocyclohexylphosphonate 1-(butylamino)-cyclohexylphosphonate de dibutyle dibutyl [1-(butylamino)cyclohexyl]= phosphonate			**P**
		$C_{18}H_{38}NO_3P$	51249-05-9	
	InChI=1/C18H38NO3P/c1-4-7-15-19-18(13-11-10-12-14-18)23(20,21-16-8-5-2)22-17-9-6-3/h19H,4-17H2,1-3H3			
E buprofezin **F** buprofézin (f)	(*Z*)-2-*tert*-butylimino-3-isopropyl-5-phenyl-1,3,5-thiadiazinan-4-one (*Z*)-2-*tert*-butylimino-3-isopropyl-5-phényl-(4*H*)1,3,5-thiadiazin-4-one (*Z*)-2-[(1,1-dimethylethyl)imino]= tetrahydro-3-(1-methylethyl)-5-phenyl-4*H*-1,3,5-thiadiazin-4-one			**IGR**
		$C_{16}H_{23}N_3OS$	953030-84-7	
	InChI=1/C16H23N3OS/c1-12(2)19-14(17-16(3,4)5)21-11-18(15(19)20)13-9-7-6-8-10-13/h6-10,12H,11H2,1-5H3/b17-14-			
E butachlor **F** butachlor (m)	*N*-butoxymethyl-2-chloro-2′,6′-diethylacetanilide *N*-(butoxyméthyl)-2-chloro-2′,6′-diéthylacétanilide *N*-(butoxymethyl)-2-chloro-*N*-(2,6-diethylphenyl)acetamide			**H**
		$C_{17}H_{26}ClNO_2$	23184-66-9	
	InChI=1/C17H26ClNO2/c1-4-7-11-21-13-19(16(20)12-18)17-14(5-2)9-8-10-15(17)6-3/h8-10H,4-7,11-13H2,1-3H3			
E carbosulfan **F** carbosulfan (m)	2,3-dihydro-2,2-dimethylbenzofuran-7-yl (dibutylaminothio)methylcarbamate (dibutylamino)thiométhylcarbamate de 2,2-diméthyl-2,3-dihydrobenzofuran-7-yle 2,3-dihydro-2,2-dimethyl-7-benzofuranyl [(dibutylamino)thio]methylcarbamate			**I** **N**
		$C_{20}H_{32}N_2O_3S$	55285-14-8	
	InChI=1/C20H32N2O3S/c1-6-8-13-22(14-9-7-2)26-21(5)19(23)24-17-12-10-11-16-15-20(3,4)25-18(16)17/h10-12H,6-9,13-15H2,1-5H3			

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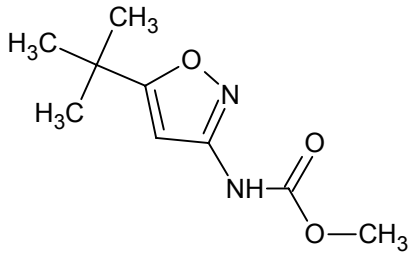
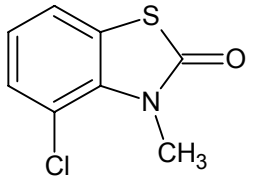
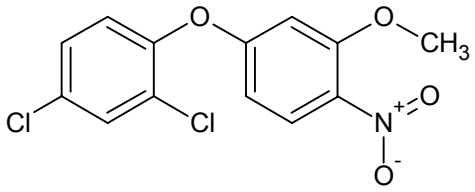
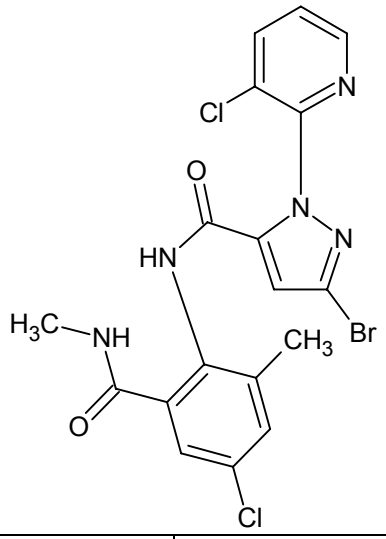
	IUPAC International Chemical Identifier (InChI™)			
E carboxazole **F** carboxazole (m)	methyl 5-*tert*-butyl-1,2-oxazol-3-ylcarbamate 5-*tert*-butylisoxazol-3-yl-carbamate de méthyle methyl [5-(1,1-dimethylethyl)-3-isoxazolyl]carbamate			**H**
		$C_9H_{14}N_2O_3$	55808-13-4	
	InChI=1/C9H14N2O3/c1-9(2,3)6-5-7(11-14-6)10-8(12)13-4/h5H,1-4H3,(H,10,11,12)/f/h10H			
E chlobenthiazone **F** chlobenthiazone (f)	4-chloro-3-methyl-1,3-benzothiazol-2(3*H*)-one 4-chloro-3-méthyl-(3*H*)-1,3-benzothiazol-2-one 4-chloro-3-methyl-2(3*H*)-benzothiazolone			**F**
		C_8H_6ClNOS	63755-05-5	
	InChI=1/C8H6ClNOS/c1-10-7-5(9)3-2-4-6(7)12-8(10)11/h2-4H,1H3			
E chlomethoxyfen **F** chlométhoxyfène (m)	5-(2,4-dichlorophenoxy)-2-nitroanisole 5-(2,4-dichlorophénoxy)-2-nitroanisole 2,4-dichloro-1-(3-methoxy-4-nitrophenoxy)benzene			**H**
		$C_{13}H_9Cl_2NO_4$	32861-85-1	
	InChI=1/C13H9Cl2NO4/c1-19-13-7-9(3-4-11(13)16(17)18)20-12-5-2-8(14)6-10(12)15/h2-7H,1H3			
E chlorantraniliprole **F** chlorantraniliprole (m)	3-bromo-4′-chloro-1-(3-chloro-2-pyridyl)-2′-methyl-6′-(methylcarbamoyl)pyrazole-5-carboxanilide 3-bromo-*N*-[4-chloro-2-méthyl-6-(méthylcarbamoyl)phényl]-1-(3-chloropyridin-2-yl)-1*H*-pyrazole-5-carboxamide 3-bromo-*N*-[4-chloro-2-methyl-6-[(methylamino)carbonyl]phenyl]-1-(3-chloro-2-pyridinyl)-1*H*-pyrazole-5-carboxamide			**I**
		$C_{18}H_{14}BrCl_2N_5O_2$	500008-45-7	
	InChI=1/C18H14BrCl2N5O2/c1-9-6-10(20)7-11(17(27)22-2)15(9)24-18(28)13-8-14(19)25-26(13)16-12(21)4-3-5-23-16/h3-8H,1-2H3,(H,22,27)(H,24,28)/f/h22,24H			

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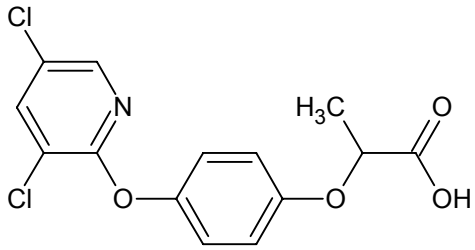
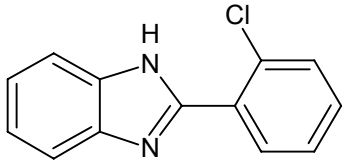
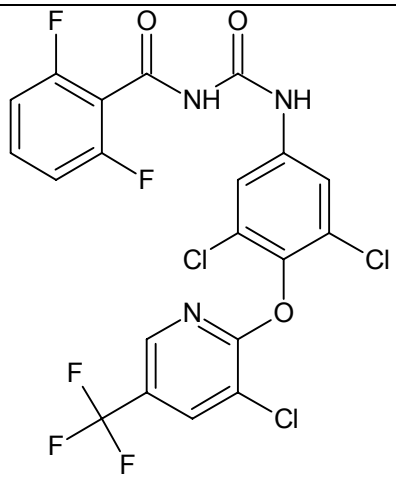
	IUPAC International Chemical Identifier (InChI™)			
E chlorazifop **F** chlorazifop (m)	(*RS*)-2-[4-(3,5-dichloro-2-pyridyloxy)phenoxy]propionic acid acide (*RS*)-2-[4-(3,5-dichloropyrid-2-yloxy)phénoxy]propionique 2-[4-[(3,5-dichloro-2-pyridinyl)oxy]phenoxy]propanoic acid	 $C_{14}H_{11}Cl_2NO_4$	60074-25-1	**H**
	InChI=1/C14H11Cl2NO4/c1-8(14(18)19)20-10-2-4-11(5-3-10)21-13-12(16)6-9(15)7-17-13/h2-8H,1H3,(H,18,19)/t8-/s3/f/h18H chlorazifop-propargyl InChI=1/C17H13Cl2NO4/c1-3-8-22-17(21)11(2)23-13-4-6-14(7-5-13)24-16-15(19)9-12(18)10-20-16/h1,4-7,9-11H,8H2,2H3/t11-/s3			
	NOTE It should be stated which ester or salt is present, for example chlorazifop-propargyl [72880-52-5]. NOTE Il convient de préciser quel est l'ester ou le sel présent, par exemple chlorazifop-propargyl [72880-52-5].			
E chlorfenazole **F** chlorfénazole (m)	2-(2-chlorophenyl)benzimidazole 2-(2-chlorophényl)benzimidazole 2-(2-chlorophenyl)-1*H*-benzimidazole	 $C_{13}H_9ClN_2$	3574-96-7	**F**
	InChI=1/C13H9ClN2/c14-10-6-2-1-5-9(10)13-15-11-7-3-4-8-12(11)16-13/h1-8H,(H,15,16)/f/h15H			
E chlorfluazuron **F** chlorfluazuron (m)	1-[3,5-dichloro-4-(3-chloro-5-trifluoromethyl-2-pyridyloxy)phenyl]-3-(2,6-difluorobenzoyl)urea 1-[3,5-dichloro-4-(3-chloro-5-trifluorométhylpyrid-2-yloxy)phényl]-3-(2,6-difluorobenzoyl)urée *N*-[[[3,5-dichloro-4-[[3-chloro-5-(trifluoromethyl)-2-pyridinyl]oxy]phenyl]amino]carbonyl]-2,6-difluorobenzamide	 $C_{20}H_9Cl_3F_5N_3O_3$	71422-67-8	**IGR**
	InChI=1/C20H9Cl3F5N3O3/c21-10-5-9(30-19(33)31-17(32)15-13(24)2-1-3-14(15)25)6-11(22)16(10)34-18-12(23)4-8(7-29-18)20(26,27)28/h1-7H,(H2,30,31,32,33)/f/h30-31H			

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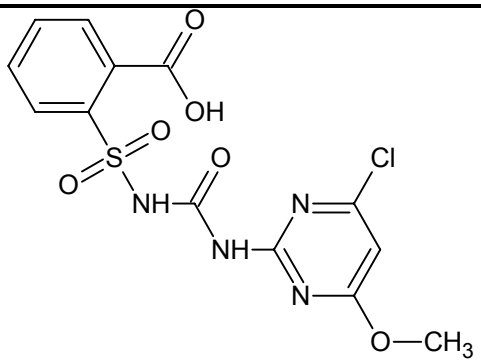
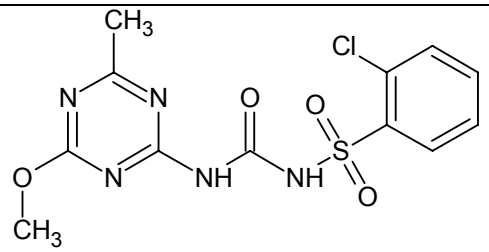
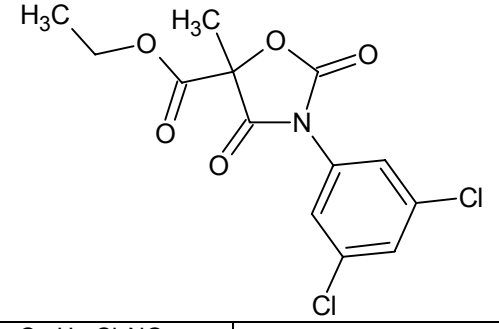
	IUPAC International Chemical Identifier (InChI™)			
E chlorimuron **F** chlorimuron (m)	2-(4-chloro-6-methoxypyrimidin-2-ylcarbamoylsulfamoyl)benzoic acid acide 2-(4-chloro-6-méthoxypyrimidin-2-yluréidosulfonyl)benzoïque 2-[[[[(4-chloro-6-methoxy-2-pyrimidinyl)amino]carbonyl]amino]=sulfonyl]benzoic acid			**H**
		$C_{13}H_{11}ClN_4O_6S$	99283-00-8	
	InChI=1/C13H11ClN4O6S/c1-24-10-6-9(14)15-12(16-10)17-13(21)18-25(22,23)8-5-3-2-4-7(8)11(19)20/h2-6H,1H3,(H,19,20)(H2,15,16,17,18,21)/f/h17-19H chlorimuron-ethyl InChI=1/C15H15ClN4O6S/c1-3-26-13(21)9-6-4-5-7-10(9)27(23,24)20-15(22)19-14-17-11(16)8-12(18-14)25-2/h4-8H,3H2,1-2H3,(H2,17,18,19,20,22)/f/h19-20H			
	NOTE It should be stated which ester or salt is present, for example chlorimuron-ethyl [90982-32-4]. NOTE Il convient de préciser quel est l'ester ou le sel présent, par exemple chlorimuron-éthyle [90982-32-4].			
E chlorsulfuron **F** chlorsulfuron (m)	1-(2-chlorophenylsulfonyl)-3-(4-methoxy-6-methyl-1,3,5-triazin-2-yl)urea 1-(2-chlorophénylsulfonyl)-3-(4-méthoxy-6-méthyl-1,3,5-triazin-2-yl)urée 2-chloro-*N*-[[(4-methoxy-6-methyl-1,3,5-triazin-2-yl)amino]carbonyl]=benzenesulfonamide			**H**
		$C_{12}H_{12}ClN_5O_4S$	64902-72-3	
	InChI=1/C12H12ClN5O4S/c1-7-14-10(17-12(15-7)22-2)16-11(19)18-23(20,21)9-6-4-3-5-8(9)13/h3-6H,1-2H3,(H2,14,15,16,17,18,19)/f/h16,18H			
E chlozolinate **F** chlozolinate (m)	ethyl (*RS*)-3-(3,5-dichlorophenyl)-5-methyl-2,4-dioxo-1,3-oxazolidine-5-carboxylate (*RS*)-3-(3,5-dichlorophényl)-5-méthyl-2,4-dioxo-1,3-oxazolidine-5-carboxylate d'éthyle ethyl 3-(3,5-dichlorophenyl)-5-methyl-2,4-dioxo-5-oxazolidinecarboxylate			**F**
		$C_{13}H_{11}Cl_2NO_5$	84332-86-5	
	InChI=1/C13H11Cl2NO5/c1-3-20-11(18)13(2)10(17)16(12(19)21-13)9-5-7(14)4-8(15)6-9/h4-6H,3H2,1-2H3/t13-/s3			

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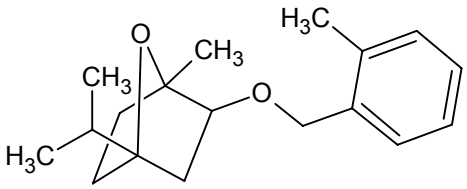
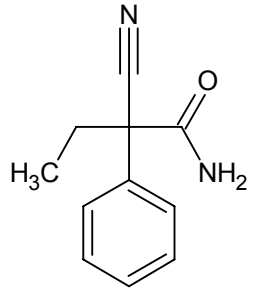
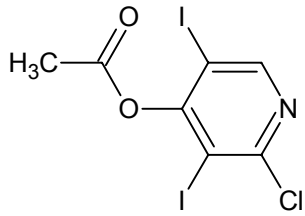
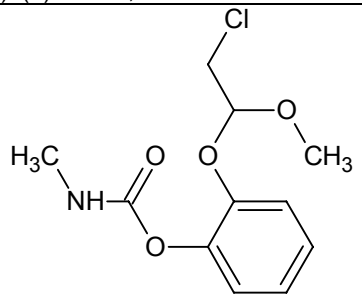
E cinmethylin **F** cinméthyline (f)	(1*RS*,2*SR*,4*SR*)-1,4-epoxy-*p*-menth-2-yl 2-methylbenzyl ether oxyde (ou éther) de (1*RS*,2*SR*,4*SR*)-1,4-époxy-*p*-menth-2-yle et de 2-méthylbenzyle (1*R*,2*S*,4*S*)-*rel*-1-methyl-4-(1-methylethyl)-2-[(2-methylphenyl)methoxy]-7-oxabicyclo[2.2.1]heptane			**H**
		$C_{18}H_{26}O_2$	87818-31-3	
	InChI=1/C18H26O2/c1-13(2)18-10-9-17(4,20-18)16(11-18)19-12-15-8-6-5-7-14(15)3/h5-8,13,16H,9-12H2,1-4H3/t16-,17-,18-/s3			
E ciobutide **F** ciobutide (m)	(*RS*)-2-cyano-2-phenylbutyramide (*RS*)-2-cyano-2-phénylbutyramide α-cyano-α-ethylbenzeneacetamide			**P**
		$C_{11}H_{12}N_2O$	80544-75-8	
	InChI=1/C11H12N2O/c1-2-11(8-12,10(13)14)9-6-4-3-5-7-9/h3-7H,2H2,1H3,(H2,13,14)/t11-/s3/f/h13H2			
E cliodinate **F** cliodinate (m)	2-chloro-3,5-diiodo-4-pyridyl acetate acétate de 2-chloro-3,5-diiodopyrid-4-yle 2-chloro-3,5-diiodo-4-pyridinyl acetate			**H**
		$C_7H_4ClI_2NO_2$	69148-12-5	
	InChI=1/C7H4ClI2NO2/c1-3(12)13-6-4(9)2-11-7(8)5(6)10/h2H,1H3			
E cloethocarb **F** cloéthocarbe (m)	2-[(*RS*)-2-chloro-1-methoxyethoxy]phenyl methylcarbamate méthylcarbamate de (*RS*)-2-(2-chloro-1-méthoxyéthoxy)phényle 2-(2-chloro-1-methoxyethoxy)phenyl methylcarbamate			**I** **M** **N**
		$C_{11}H_{14}ClNO_4$	51487-69-5	
	InChI=1/C11H14ClNO4/c1-13-11(14)17-9-6-4-3-5-8(9)16-10(7-12)15-2/h3-6,10H,7H2,1-2H3,(H,13,14)/t10-/s3/f/h13H			

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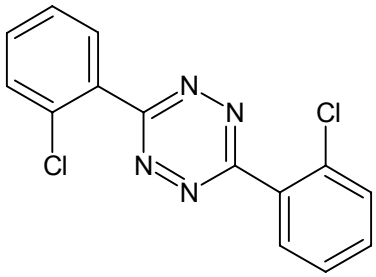
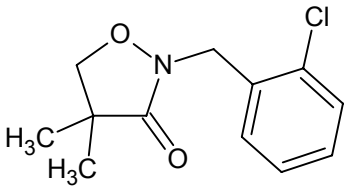
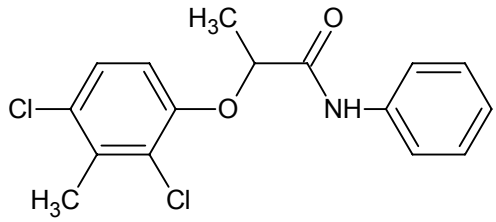
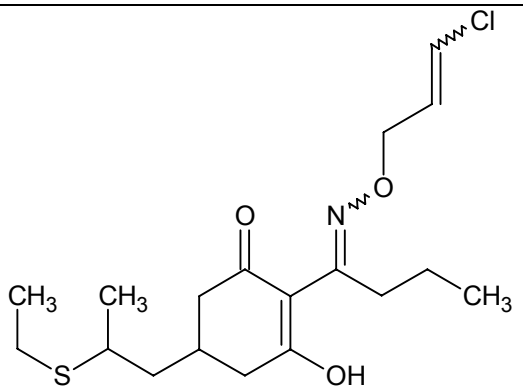
E clofentezine **F** clofentézine (f)	3,6-bis(2-chlorophenyl)-1,2,4,5-tetrazine 3,6-bis(2-chlorophényl)-1,2,4,5-tétrazine 3,6-bis(2-chlorophenyl)-1,2,4,5-tetrazine			**A**
		$C_{14}H_8Cl_2N_4$	74115-24-5	
	InChI=1/C14H8Cl2N4/c15-11-7-3-1-5-9(11)13-17-19-14(20-18-13)10-6-2-4-8-12(10)16/h1-8H			
E clomazone **F** clomazone (f)	2-(2-chlorobenzyl)-4,4-dimethyl-1,2-oxazolidin-3-one 2-(2-chlorobenzyl)-4,4-diméthyl-1,2-oxazolidin-3-one 2-[(2-chlorophenyl)methyl]-4,4-dimethyl-3-isoxazolidinone			**H**
		$C_{12}H_{14}ClNO_2$	81777-89-1	
	InChI=1/C12H14ClNO2/c1-12(2)8-16-14(11(12)15)7-9-5-3-4-6-10(9)13/h3-6H,7-8H2,1-2H3			
E clomeprop **F** cloméprop (m)	(*RS*)-2-(2,4-dichloro-*m*-tolyloxy)propionanilide (*RS*)-2-(2,4-dichloro-3-méthylphénoxy)propionanilide 2-(2,4-dichloro-3-methylphenoxy)-*N*-phenylpropanamide			**H**
		$C_{16}H_{15}Cl_2NO_2$	84496-56-0	
	InChI=1/C16H15Cl2NO2/c1-10-13(17)8-9-14(15(10)18)21-11(2)16(20)19-12-6-4-3-5-7-12/h3-9,11H,1-2H3,(H,19,20)/t11-/s3/f/h19H			
E cloproxydim **F** cloproxydime (f)	(5*RS*)-2-{(*EZ*)-1-[(2*EZ*)-3-chloroallyloxyimino]butyl}-5-[(2*RS*)-2-ethylthiopropyl]-3-hydroxycyclohex-2-en-1-one (5*RS*)-2-{(*EZ*)-1-[(2*EZ*)-3-chloroallyloxyimino]butyl}-5-[(2*RS*)-2-éthylthiopropyl]-3-hydroxycyclohex-2-èn-1-one 2-[1-[[(3-chloro-2-propenyl)oxy]imino]butyl]-5-[2-(ethylthio)propyl]-3-hydroxy-2-cyclohexen-1-one			**H**
		$C_{18}H_{28}ClNO_3S$	95480-33-4	
	InChI=1/C18H28ClNO3S/c1-4-7-15(20-23-9-6-8-19)18-16(21)11-14(12-17(18)22)10-13(3)24-5-2/h6,8,13-14,21H,4-5,7,9-12H2,1-3H3/b8-6?,20-15?/t13?,14?			

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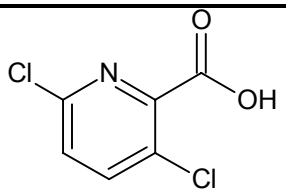
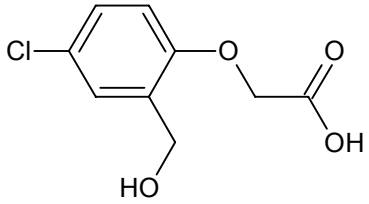
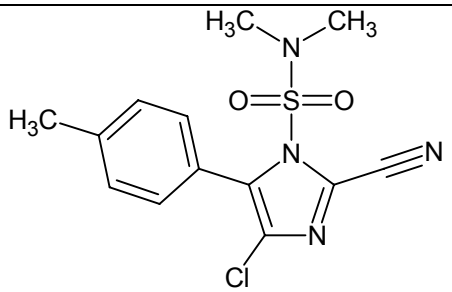
E clopyralid **F** clopyralid (m)	3,6-dichloropyridine-2-carboxylic acid acide 3,6-dichloropicolique ou 3,6-dichloropicolinique 3,6-dichloro-2-pyridinecarboxylic acid			**H**
		$C_6H_3Cl_2NO_2$	1702-17-6	
	InChI=1/C6H3Cl2NO2/c7-3-1-2-4(8)9-5(3)6(10)11/h1-2H,(H,10,11)/f/h10H clopyralid-olamine InChI=1/C6H3Cl2NO2.C2H7NO/c7-3-1-2-4(8)9-5(3)6(10)11;3-1-2-4/h1-2H,(H,10,11);4H,1-3H2/f/h10H; clopyralid-methyl InChI=1/C7H5Cl2NO2/c1-12-7(11)6-4(8)2-3-5(9)10-6/h2-3H,1H3			
	NOTE 1 It should be stated which ester or salt is present, for example clopyralid-methyl [1532-24-7] or clopyralid-olamine [57754-85-5]. NOTE 1 Il convient de préciser quel est l'ester ou le sel présent, par exemple clopyralid-méthyle [1532-24-7] ou clopyralid-olamine [57754-85-5]. NOTE 2 The name “clopyralid” is not acceptable for use in Canada, Finland and France, where the shortened chemical name “3,6-dichloropicolinic acid” is considered suitable. NOTE 2 Le nom «clopyralid» n'est pas acceptable pour l'emploi au Canada, en Finlande et en France, où le nom chimique acide 3,6-dichloropicolinique est considéré approprié.			
E cloxyfonac **F** cloxyfonac (m)	4-chloro-α-hydroxy-*o*-tolyloxyacetic acid acide 4-chloro-2-hydroxyméthylphénoxyacétique [4-chloro-2-(hydroxymethyl)phenoxy]acetic acid			**P**
		$C_9H_9ClO_4$	6386-63-6	
	InChI=1/C9H9ClO4/c10-7-1-2-8(6(3-7)4-11)14-5-9(12)13/h1-3,11H,4-5H2,(H,12,13)/f/h12H cloxyfonac-sodium InChI=1/C9H9ClO4.Na/c10-7-1-2-8(6(3-7)4-11)14-5-9(12)13;/h1-3,11H,4-5H2,(H,12,13);/q;+1/p-1/fC9H8ClO4.Na/q-1;m			
	NOTE It should be stated which ester or salt is present, for example cloxyfonac-sodium [32791-87-0]. NOTE Il convient de préciser quel est l'ester ou le sel présent, par exemple cloxyfonac-sodium [32791-87-0].			
E cyazofamid **F** cyazofamide (m)	4-chloro-2-cyano-*N,N*-dimethyl-5-*p*-tolylimidazole-1-sulfonamide *N,N*-diméthyl-(4-chloro-2-cyano-5-*p*-tolylimidazole)-1-sulfonamide 4-chloro-2-cyano-*N,N*-dimethyl-5-(4-methylphenyl)-1*H*-imidazole-1-sulfonamide			**F**
		$C_{13}H_{13}ClN_4O_2S$	120116-88-3	
	InChI=1/C13H13ClN4O2S/c1-9-4-6-10(7-5-9)12-13(14)16-11(8-15)18(12)21(19,20)17(2)3/h4-7H,1-3H3			

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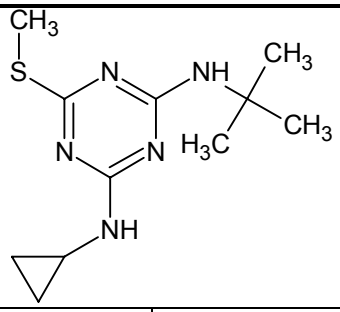
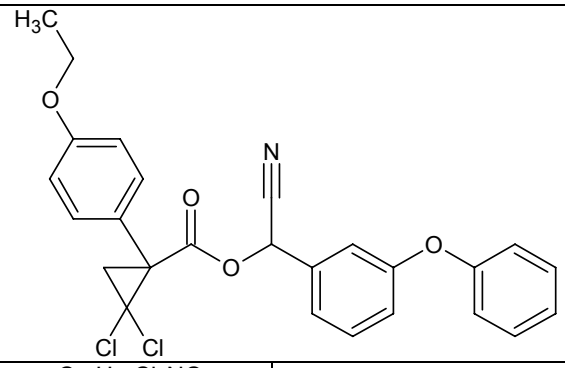
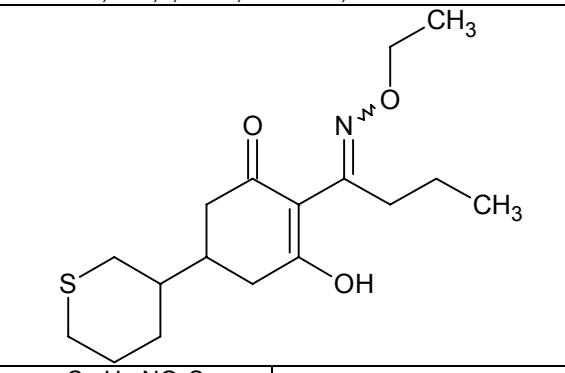
E cybutryne **F** cybutryne (m)	N^2-*tert*-butyl-N^4-cyclopropyl-6-methylthio-1,3,5-triazine-2,4-diamine *N*-*tert*-butyl-*N'*-cyclopropyl-6-méthylthio-1,3,5-triazine-2,4-diamine *N*-cyclopropyl-*N'*-(1,1-dimethylethyl)-6-(methylthio)-1,3,5-triazine-2,4-diamine			**AL**
		$C_{11}H_{19}N_5S$	28159-98-0	
	InChI=1/C11H19N5S/c1-11(2,3)16-9-13-8(12-7-5-6-7)14-10(15-9)17-4/h7H,5-6H2,1-4H3,(H2,12,13,14,15,16)/f/h12,16			
E cycloprothrin **F** cycloprothrine (f)	(*RS*)-α-cyano-3-phenoxybenzyl (*RS*)-2,2-dichloro-1-(4-ethoxyphenyl)=cyclopropanecarboxylate (*RS*)-2,2-dichloro-1-(4-éthoxyphényl)-cyclopropanecarboxylate de (*RS*)-1-cyano-1-(3-phénoxyphényl)méthyle cyano(3-phenoxyphenyl)methyl 2,2-dichloro-1-(4-ethoxyphenyl)=cyclopropanecarboxylate			**I**
		$C_{26}H_{21}Cl_2NO_4$	63935-38-6	
	InChI=1/C26H21Cl2NO4/c1-2-31-20-13-11-19(12-14-20)25(17-26(25,27)28)24(30)33-23(16-29)18-7-6-10-22(15-18)32-21-8-4-3-5-9-21/h3-15,23H,2,17H2,1H3/t23?,25?			
E cycloxydim **F** cycloxydime (f)	(5*RS*)-2-[(*EZ*)-1-(ethoxyimino)butyl]-3-hydroxy-5-[(3*RS*)-thian-3-yl]cyclohex-2-en-1-one (5*RS*)-2-[(*EZ*)-1-(éthoxyimino)butyl]-3-hydroxy-5-[(3*RS*)-thian-3-yl]cyclohex-2-èn-1-one 2-[1-(ethoxyimino)butyl]-3-hydroxy-5-(tetrahydro-2*H*-thiopyran-3-yl)-2-cyclohexen-1-one			**H**
		$C_{17}H_{27}NO_3S$	101205-02-1	
	InChI=1/C17H27NO3S/c1-3-6-14(18-21-4-2)17-15(19)9-13(10-16(17)20)12-7-5-8-22-11-12/h12-13,19H,3-11H2,1-2H3/b18-14?/t12?,13?			

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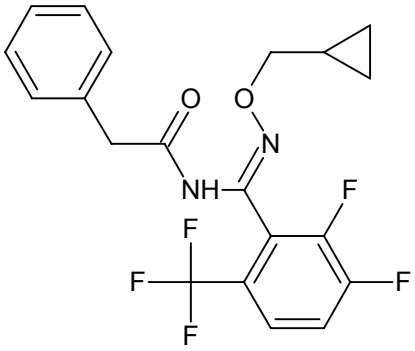
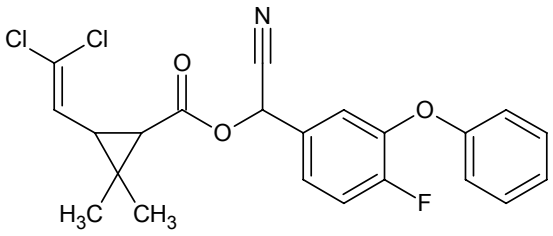
	IUPAC International Chemical Identifier (InChI™)			
E cyflufenamid **F** cyfluf énamide (m)	(*Z*)-*N*-[α-(cyclopropylmethoxyimino)-2,3-difluoro-6-(trifluoromethyl)benzyl]-2-phenylacetamide (*Z*)-*N*-[(cyclopropyl(méthoxy)imino)-(2,3-difluoro-6-trifluorométhylphényl)-méthyl]-2-phénylacétamide [*N*(*Z*)]-*N*-[[(cyclopropylmethoxy)=amino][2,3-difluoro-6-(trifluoromethyl)phenyl]methylene]=benzeneacetamide			**F**
		$C_{20}H_{17}F_5N_2O_2$	180409-60-3	
	InChI=1/C20H17F5N2O2/c21-15-9-8-14(20(23,24)25)17(18(15)22)19(27-29-11-13-6-7-13)26-16(28)10-12-4-2-1-3-5-12/h1-5,8-9,13H,6-7,10-11H2,(H,26,27,28)/f/h26H/b27-19- CAS-preferred tautomer InChI=1/C20H17F5N2O2/c21-15-9-8-14(20(23,24)25)17(18(15)22)19(27-29-11-13-6-7-13)26-16(28)10-12-4-2-1-3-5-12/h1-5,8-9,13H,6-7,10-11H2,(H,26,27,28)/f/h27H/b26-19-			
	NOTE The CAS name and Registry Number are for the tautomer that is preferred under the CAS nomenclature rules. NOTE Le nom et le numéro de registre de CAS sont attribués au tautomère préféré conformément aux règles de nomenclature du CAS.			
E cyfluthrin **F** cyfluthrine (f)	(*RS*)-α-cyano-4-fluoro-3-phenoxybenzyl (1*RS,*3*RS*;1*RS,*3*SR*)-3-(2,2-dichlorovinyl)-2,2-dimethylcyclopropanecarboxylate Alternative Rothamsted-style stereodescriptors: (*RS*)-α-cyano-4-fluoro-3-phenoxybenzyl (1*RS*)-*cis-trans*-3-(2,2-dichlorovinyl)-2,2-dimethylcyclopropanecarboxylate (1*RS,*3*RS*;1*RS,*3*SR*)-3-(2,2-dichlorovinyl)-2,2-diméthylcyclopropanecarboxylate de (*RS*)-cyano-4-fluoro-3-phénoxyphénylméthyle cyano(4-fluoro-3-phenoxyphenyl)methyl 3-(2,2-dichloroethenyl)-2,2-dimethylcyclopropanecarboxylate			**I**
		$C_{22}H_{18}Cl_2FNO_3$	68359-37-5	
	InChI=1/C22H18Cl2FNO3/c1-22(2)15(11-19(23)24)20(22)21(27)29-18(12-26)13-8-9-16(25)17(10-13)28-14-6-4-3-5-7-14/h3-11,15,18,20H,1-2H3/t15?,18?,20?			

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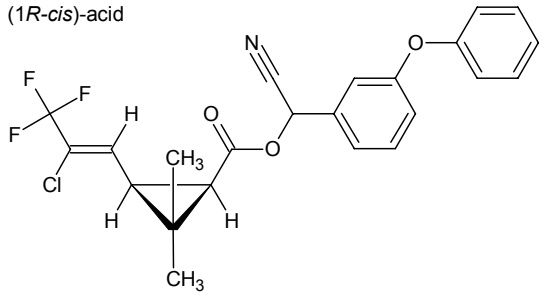
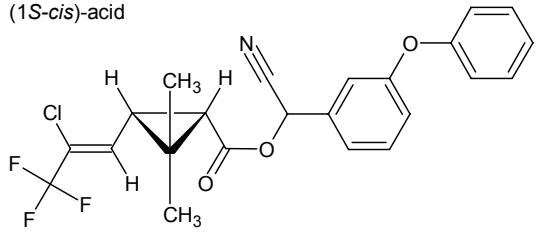
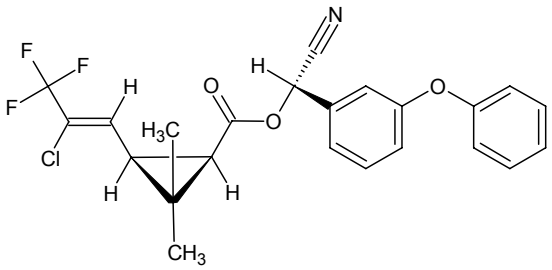
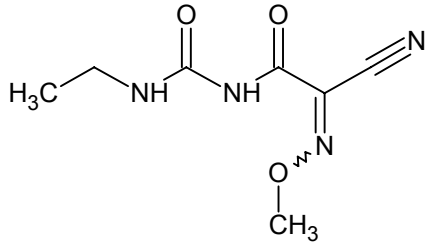
E cyhalothrin **F** cyhalothrine (f)	(*RS*)-α-cyano-3-phenoxybenzyl (1*RS*,3*RS*)-3-[(*Z*)-2-chloro-3,3,3-trifluoropropenyl]-2,2-dimethylcyclopropanecarboxylate Alternative Rothamsted-style stereodescriptors: (*RS*)-α-cyano-3-phenoxybenzyl (1*RS*)-*cis*-3-[(*Z*)-2-chloro-3,3,3-trifluoropropenyl]-2,2-dimethylcyclopropanecarboxylate	(1*R-cis*)-acid 		**A** **I**
	(1*RS*,3*RS*)-3-[(*Z*)-2-chloro-3,3,3-trifluoroprop-1-ényl]-2,2-diméthylcyclopropanecarboxylate de (*RS*)-cyano-3-phénoxyphénylméthyle	(1*S-cis*)-acid 		
	cyano(3-phenoxyphenyl)methyl (1*R*,3*R*)-*rel*-3-[(1*Z*)-2-chloro-3,3,3-trifluoro-1-propenyl]-2,2-dimethylcyclopropanecarboxylate	$C_{23}H_{19}ClF_3NO_3$	68085-85-8	
	InChI=1/C23H19ClF3NO3/c1-22(2)17(12-19(24)23(25,26)27)20(22)21(29)31-18(13-28)14-7-6-10-16(11-14)30-15-8-4-3-5-9-15/h3-12,17-18,20H,1-2H3/b19-12-/t17-,18?,20-/s3			
E gamma-cyhalothrin **F** gamma-cyhalothrine (f)	(*S*)-α-cyano-3-phenoxybenzyl (1*R*,3*R*)-3-[(*Z*)-2-chloro-3,3,3-trifluoropropenyl]-2,2-dimethylcyclopropanecarboxylate Alternative Rothamsted-style stereodescriptors: (*S*)-α-cyano-3-phenoxybenzyl (1*R*)-*cis*-3-[(*Z*)-2-chloro-3,3,3-trifluoropropenyl]-2,2-dimethylcyclopropanecarboxylate			**I**
	(1*R*,3*R*)-3-[(*Z*)-2-chloro-3,3,3-trifluoropropényl]-2,2-diméthylcyclopropanecarboxylate de (*S*)-cyano-(3-phénoxyphényl)-méthyle			
	(*S*)-cyano(3-phenoxyphenyl)methyl (1*R*,3*R*)-3-[(1*Z*)-2-chloro-3,3,3-trifluoro-1-propenyl]-2,2-dimethylcyclopropanecarboxylate	$C_{23}H_{19}ClF_3NO_3$	76703-62-3	
	InChI=1/C23H19ClF3NO3/c1-22(2)17(12-19(24)23(25,26)27)20(22)21(29)31-18(13-28)14-7-6-10-16(11-14)30-15-8-4-3-5-9-15/h3-12,17-18,20H,1-2H3/b19-12-/t17-,18+,20-/m0/s1			
E cymoxanil **F** cymoxanil (m)	1-[(*EZ*)-2-cyano-2-methoxyiminoacetyl]-3-ethylurea			**F**
	1-[(*EZ*)-cyano-2-méthoxyiminoacétyl]-3-éthylurée			
	2-cyano-*N*-[(ethylamino)carbonyl]-2-(methoxyimino)acetamide	$C_7H_{10}N_4O_3$	57966-95-7	
	InChI=1/C7H10N4O3/c1-3-9-7(13)10-6(12)5(4-8)11-14-2/h3H2,1-2H3,(H2,9,10,12,13)/b11-5?/f/h9-10H			

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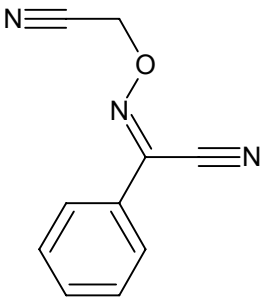
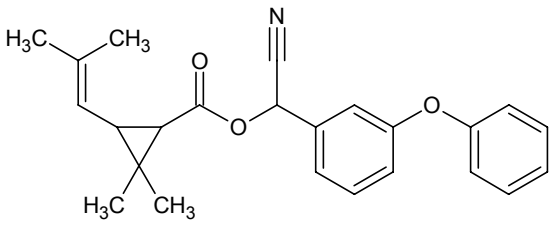
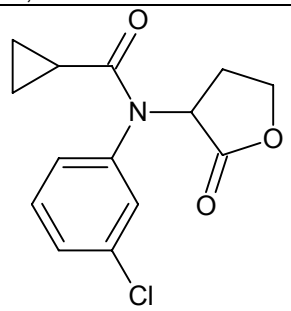
	IUPAC International Chemical Identifier (InChI™)			
E cyometrinil **F** cyométrinil (m)	(*Z*)-cyanomethoxyimino=(phenyl)acetonitrile (*Z*)-cyanométhoxyimino-(phényl)acétonitrile (α*Z*)-α-[(cyanomethoxy)imino]=benzeneacetonitrile			**S**
		$C_{10}H_7N_3O$	78370-21-5	
	InChI=1/C10H7N3O/c11-6-7-14-13-10(8-12)9-4-2-1-3-5-9/h1-5H,7H2/b13-10+			
E cyphenothrin **F** cyphénothrine (f)	(*RS*)-α-cyano-3-phenoxybenzyl (1*RS*,3*RS*;1*RS*,3*SR*)-2,2-dimethyl-3-(2-methylprop-1-enyl)cyclopropanecarboxylate Alternative Rothamsted-style stereodescriptors: (*RS*)-α-cyano-3-phenoxybenzyl (1*RS*)-*cis-trans*-2,2-dimethyl-3-(2-methylprop-1-enyl)cyclopropanecarboxylate (1*RS*,3*RS*;1*RS*,3*SR*)-2,2-diméthyl-3-(2-méthylprop-1-ènyl)-cyclopropanecarboxylate de (*RS*)-cyano-(3-phénoxyphényl)méthyle cyano(3-phenoxyphenyl)methyl 2,2-dimethyl-3-(2-methyl-1-propenyl)cyclopropanecarboxylate			**I**
		$C_{24}H_{25}NO_3$	39515-40-7	
	InChI=1/C24H25NO3/c1-16(2)13-20-22(24(20,3)4)23(26)28-21(15-25)17-9-8-12-19(14-17)27-18-10-6-5-7-11-18/h5-14,20-22H,1-4H3/t20?,21?,22?			
E cyprofuram **F** cyprofuram (m)	(*RS*)-α-[*N*-(3-chlorophenyl)=cyclopropanecarboxamido]-γ-butyrolactone (*RS*)-*N*-(3-chlorophényl)-*N*-(γ-butyrolacton-3-yl)-cyclopropanecarboxamide *N*-(3-chlorophenyl)-*N*-(tetrahydro-2-oxo-3-furanyl)cyclopropanecarboxamide			**F**
		$C_{14}H_{14}ClNO_3$	69581-33-5	
	InChI=1/C14H14ClNO3/c15-10-2-1-3-11(8-10)16(13(17)9-4-5-9)12-6-7-19-14(12)18/h1-3,8-9,12H,4-7H2/t12-/s3			

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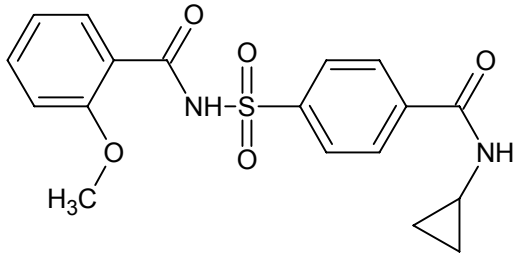
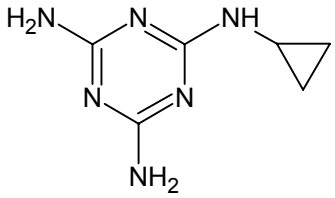
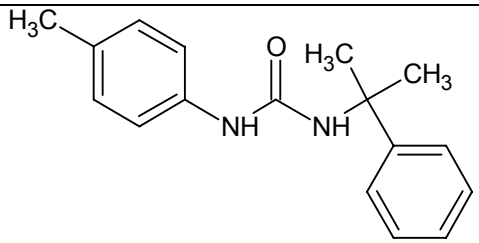
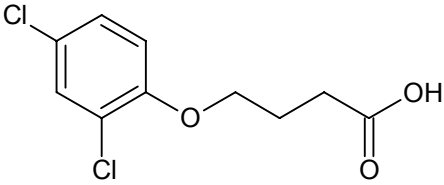
E cyprosulfamide **F** cyprosulfamide (m)	*N*-[4-(cyclopropylcarbamoyl)= phenylsulfonyl]-*o*-anisamide *N*-[4-(cyclopropylcarbamoyl)= phénylsulfonyl]-*o*-anisamide *N*-[[4-[(cyclopropylamino)carbonyl]= phenyl]sulfonyl]-2-methoxybenzamide			**P** **S**
		$C_{18}H_{18}N_2O_5S$	221667-31-8	
	InChI=1/C18H18N2O5S/c1-25-16-5-3-2-4-15(16)18(22)20-26(23,24)14-10-6-12(7-11-14)17(21)19-13-8-9-13/h2-7,10-11,13H,8-9H2,1H3,(H,19,21)(H,20,22)/f/h19-20H			
E cyromazine **F** cyromazine (f)	*N*-cyclopropyl-1,3,5-triazine-2,4,6-triamine *N*-cyclopropyl-1,3,5-triazine-2,4,6-triamine *N*-cyclopropyl-1,3,5-triazine-2,4,6-triamine			**A** **IGR**
		$C_6H_{10}N_6$	66215-27-8	
	InChI=1/C6H10N6/c7-4-10-5(8)12-6(11-4)9-3-1-2-3/h3H,1-2H2,(H5,7,8,9,10,11,12)/f/h9H,7-8H2			
E daimuron **F** daimuron (m)	1-(1-methyl-1-phenylethyl)-3-*p*-tolylurea 1-(2-phénylprop-2-yl)-3-*p*-tolylurée *N*-(4-methylphenyl)-*N*′-(1-methyl-1-phenylethyl)urea			**H**
		$C_{17}H_{20}N_2O$	42609-52-9	
	InChI=1/C17H20N2O/c1-13-9-11-15(12-10-13)18-16(20)19-17(2,3)14-7-5-4-6-8-14/h4-12H,1-3H3,(H2,18,19,20)/f/h18-19H			
E 2,4-DB **F** 2,4-DB (m)	4-(2,4-dichlorophenoxy)butyric acid acide 4-(2,4-dichlorophénoxy)butyrique 4-(2,4-dichlorophenoxy)butanoic acid			**H** **P**
		$C_{10}H_{10}Cl_2O_3$	94-82-6	
	InChI=1/C10H10Cl2O3/c11-7-3-4-9(8(12)6-7)15-5-1-2-10(13)14/h3-4,6H,1-2,5H2,(H,13,14)/f/h13H 2,4-DB-sodium InChI=1/C10H10Cl2O3.Na/c11-7-3-4-9(8(12)6-7)15-5-1-2-10(13)14;/h3-4,6H,1-2,5H2,(H,13,14);/q;+1/p-1/fC10H9Cl2O3.Na/q-1;m			
	NOTE It should be stated which ester or salt is present, for example 2,4-DB-isoctyl [1320-15-6] or 2,4-DB-sodium [10433-59-7]. NOTE Il convient de préciser quel est l'ester ou le sel présent, par exemple 2,4-DB-isoctyl [1320-15-6] ou 2,4-DB-sodium [10433-59-7].			

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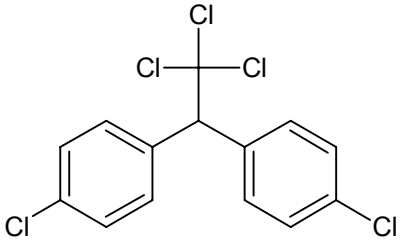
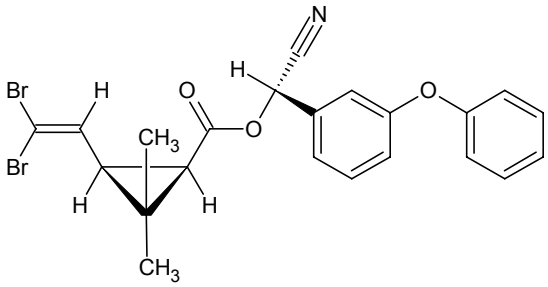
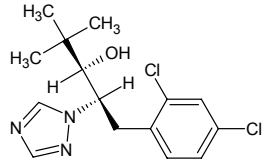
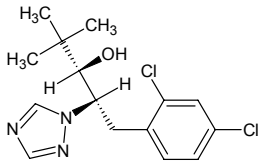
E *pp′*-DDT **F** *pp′*-DDT (m)	1,1,1-trichloro-2,2-bis(4-chlorophenyl)ethane 1,1,1-trichloro-2,2-bis(4-chlorophényl)éthane 1,1′-(2,2,2-trichloroethylidene)bis[4-chlorobenzene]			**I**
		$C_{14}H_9Cl_5$	50-29-3	
	InChI=1/C14H9Cl5/c15-11-5-1-9(2-6-11)13(14(17,18)19)10-3-7-12(16)8-4-10/h1-8,13H			
	NOTE *pp′*-DDT is the common name for the pure compound; the technical product has the ISO common name DDT. NOTE *pp′*-DDT est le nom commun du composé pur; le produit technique est appelé par son nom commun ISO «DDT».			
E deltamethrin **F** deltaméthrine (f)	(*S*)-α-cyano-3-phenoxybenzyl (1*R*,3*R*)-3-(2,2-dibromovinyl)-2,2-dimethylcyclopropanecarboxylate Alternative Rothamsted-style stereodescriptors: (*S*)-α-cyano-3-phenoxybenzyl (1*R*)-*cis*-3-(2,2-dibromovinyl)-2,2-dimethylcyclopropanecarboxylate (1*R*,3*R*)-3-(2,2-dibromovinyl)-2,2-diméthylcyclopropanecarboxylate de (*S*)-cyano-(3-phénoxyphényl)méthyle (*S*)-cyano(3-phenoxyphenyl)methyl (1*R*,3*R*)-3-(2,2-dibromoethenyl)-2,2-dimethylcyclopropanecarboxylate			**I**
		$C_{22}H_{19}Br_2NO_3$	52918-63-5	
	InChI=1/C22H19Br2NO3/c1-22(2)17(12-19(23)24)20(22)21(26)28-18(13-25)14-7-6-10-16(11-14)27-15-8-4-3-5-9-15/h3-12,17-18,20H,1-2H3/t17-,18+,20-/m0/s1			
E diclobutrazol **F** diclobutrazol (m)	(2*RS*,3*RS*)-1-(2,4-dichlorophenyl)-4,4-dimethyl-2-(1*H*-1,2,4-triazol-1-yl)pentan-3-ol (2*RS*,3*RS*)-1-(2,4-dichlorophényl)-4,4-diméthyl-2-(1*H*-1,2,4-triazol-1-yl) pentan-3-ol (α*R*,β*R*)-*rel*-β-[(2,4-dichlorophenyl)methyl]-α-(1,1-dimethylethyl)-1*H*-1,2,4-triazole-1-ethanol	 (2*R*,3*R*)-isomer  (2*S*,3*S*)-isomer		**F**
		$C_{15}H_{19}Cl_2N_3O$	75736-33-3	
	InChI=1/C15H19Cl2N3O/c1-15(2,3)14(21)13(20-9-18-8-19-20)6-10-4-5-11(16)7-12(10)17/h4-5,7-9,13-14,21H,6H2,1-3H3/t13-,14+/s3			

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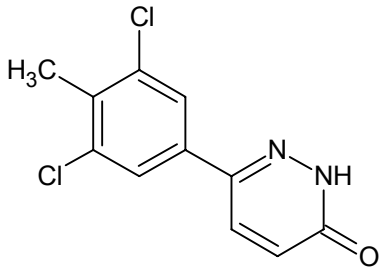
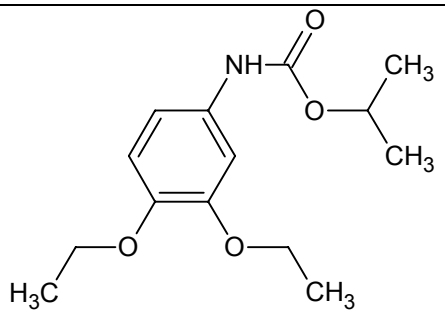
E diclomezine **F** diclomézine (f)	6-(3,5-dichlorophenyl-*p*-tolyl)pyridazin-3(2*H*)-one 6-(3,5-dichlorophényl-4-méthylphényl)-(2*H*)-pyridazin-3-one 6-(3,5-dichloro-4-methylphenyl)-3(2*H*)-pyridazinone			**F**
		$C_{11}H_8Cl_2N_2O$	62865-36-5	
	InChI=1/C11H8Cl2N2O/c1-6-8(12)4-7(5-9(6)13)10-2-3-11(16)15-14-10/h2-5H,1H3,(H,15,16)/f/h15H diclomezine-sodium InChI=1/C11H8Cl2N2O.Na/c1-6-8(12)4-7(5-9(6)13)10-2-3-11(16)15-14-10;/h2-5H,1H3,(H,15,16);/q;+1/p-1/fC11H7Cl2N2O.Na/q-1;m			
	NOTE It should be stated which salt is present, for example diclomezine-sodium [62902-57-2]. NOTE Il convient de préciser quel est le sel présent, par exemple diclomézine-sodium [62902-57-2].			
E diethofencarb **F** diéthofencarb (m)	isopropyl 3,4-diethoxycarbanilate 3,4-diéthoxycarbanilate d'isopropyle 1-methylethyl (3,4-diethoxyphenyl)carbamate			**F**
		$C_{14}H_{21}NO_4$	87130-20-9	
	InChI=1/C14H21NO4/c1-5-17-12-8-7-11(9-13(12)18-6-2)15-14(16)19-10(3)4/h7-10H,5-6H2,1-4H3,(H,15,16)/f/h15H			

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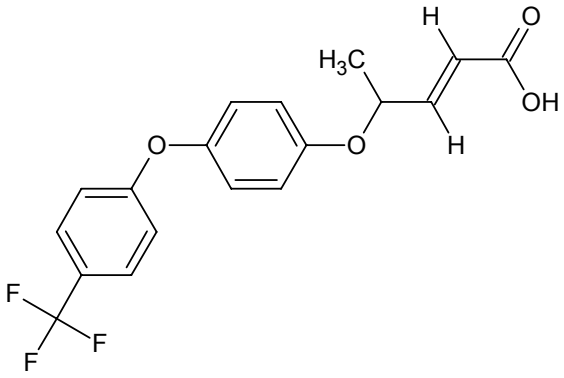
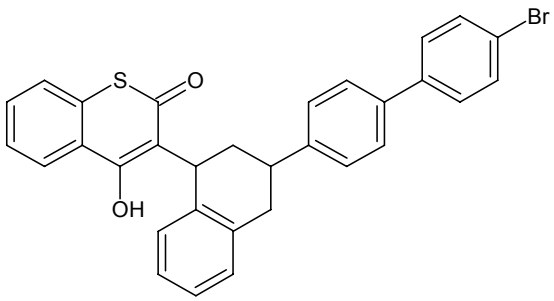
E difenopenten **F** difénopentène (m)	(*E*)-(*RS*)-4-[4-(α,α,α-trifluoro-*p*-tolyloxy)phenoxy]pent-2-enoic acid acide (*E*)-(*RS*)-4-[4-(4-trifluorométhyl)phénoxy]phénoxypent-2-ènoïque (2*E*)-4-[4-[4-(trifluoromethyl)phenoxy]phenoxy]-2-pentenoic acid	 $C_{18}H_{15}F_3O_4$	81416-44-6	**H**
	InChI=1/C18H15F3O4/c1-12(2-11-17(22)23)24-14-7-9-16(10-8-14)25-15-5-3-13(4-6-15)18(19,20)21/h2-12H,1H3,(H,22,23)/b11-2+/t12-/s3/f/h22H difenopenten-ethyl InChI=1/C20H19F3O4/c1-3-25-19(24)13-4-14(2)26-16-9-11-18(12-10-16)27-17-7-5-15(6-8-17)20(21,22)23/h4-14H,3H2,1-2H3/b13-4+/t14-/s3			
	NOTE It should be stated which ester or salt is present, for example difenopenten-ethyl [71101-05-8]. NOTE Il convient de préciser quel est l'ester ou le sel présent, par exemple difénopentène-éthyle [71101-05-8].			
E difethialone **F** difféthialone (f)	3-[(1*RS*,3*RS*;1*RS*,3*SR*)-3-(4′-bromobiphenyl-4-yl)-1,2,3,4-tetrahydro-1-naphthyl]-4-hydroxy-1-benzothiin-2-one where the ratios of the racemates (1*RS*,3*RS*) to (1*RS*,3*SR*) lie within the ranges 0–15 to 85–100 respectively 3-[(1*RS*,3*RS*;1*RS*,3*SR*)-3-(4′-bromobiphényl-4-yl)-1,2,3,4-tétrahydronapht-1-yl]-4-hydroxy-1-benzothiin-2-one Le rapport du racémate (1*RS*,3*RS*) au racémate (1*RS*,3*SR*) est situé entre 0–15 et 85–100 respectivement 3-[3-(4′-bromo[1,1′-biphenyl]-4-yl)-1,2,3,4-tetrahydro-1-naphthalenyl]-4-hydroxy-2*H*-1-benzothiopyran-2-one	 $C_{31}H_{23}BrO_2S$	104653-34-1	**R**
	InChI=1/C31H23BrO2S/c32-24-15-13-20(14-16-24)19-9-11-21(12-10-19)23-17-22-5-1-2-6-25(22)27(18-23)29-30(33)26-7-3-4-8-28(26)35-31(29)34/h1-16,23,27,33H,17-18H2/t23?,27?			
	NOTE The ratio of the racemic mixtures (1*RS*,3*RS*) to (1*RS*,3*SR*) should be stated. NOTE Il convient de préciser le rapport du racémate (1*RS*,3*RS*) au racémate (1*RS*,3*SR*).			

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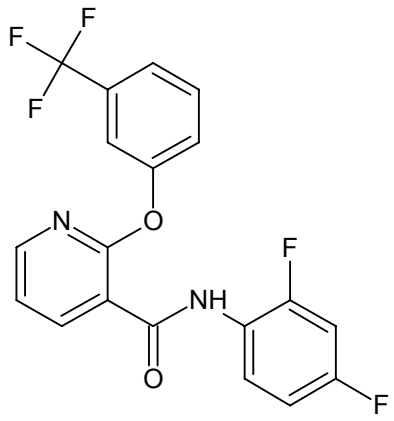
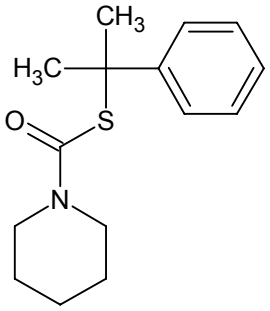
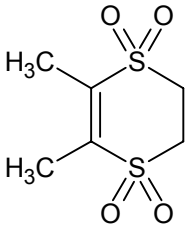
E diflufenican **F** diflufénican (m)	2′,4′-difluoro-2-(α,α,α-trifluoro-*m*-tolyloxy)nicotinanilide 2′,4′-difluoro-2-(3-trifluorométhylphénoxy)nicotinanilide *N*-(2,4-difluorophenyl)-2-[3-(trifluoromethyl)phenoxy]-3-pyridinecarboxamide			**H**
		$C_{19}H_{11}F_5N_2O_2$	83164-33-4	
	InChI=1/C19H11F5N2O2/c20-12-6-7-16(15(21)10-12)26-17(27)14-5-2-8-25-18(14)28-13-4-1-3-11(9-13)19(22,23)24/h1-10H,(H,26,27)/f/h26H			
	NOTE The name “diflufenican” is not acceptable for use in France, where “diflufénicanil” has been adopted as the common name. NOTE Le nom «diflufénican» n'est pas acceptable pour l'emploi en France, où «diflufénicanil» a été adopté comme nom commun.			
E dimepiperate **F** dimépipérate (m)	*S*-1-methyl-1-phenylethyl piperidine-1-carbothioate pipéridine-1-thiocarboxylate de *S*-2-phénylprop-2-yle *S*-(1-methyl-1-phenylethyl) 1-piperidinecarbothioate			**H**
		$C_{15}H_{21}NOS$	61432-55-1	
	InChI=1/C15H21NOS/c1-15(2,13-9-5-3-6-10-13)18-14(17)16-11-7-4-8-12-16/h3,5-6,9-10H,4,7-8,11-12H2,1-2H3			
E dimethipin **F** diméthipin (m)	2,3-dihydro-5,6-dimethyl-1,4-dithiine 1,1,4,4-tetraoxide 1,1,4,4-tétraoxyde de 2,3-dihydro-5,6-diméthyl-1,4-dithiine 2,3-dihydro-5,6-dimethyl-1,4-dithiin 1,1,4,4-tetraoxide			**P**
		$C_6H_{10}O_4S_2$	55290-64-7	
	InChI=1/C6H10O4S2/c1-5-6(2)12(9,10)4-3-11(5,7)8/h3-4H2,1-2H3			
	NOTE The name “dimethipin” is not acceptable for use in Japan because of the risk of confusion with the trade name “Dimetapin”. NOTE Le nom «diméthipin» n'est pas acceptable pour l'emploi au Japon, car il entre en conflit avec le nom commercial «Dimetapin».			

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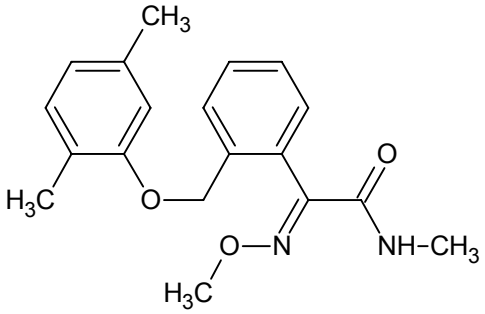
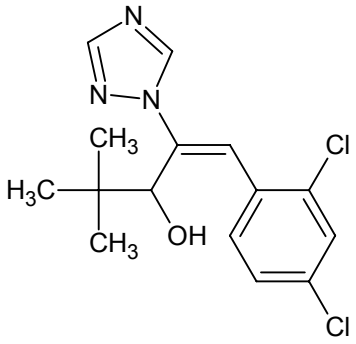
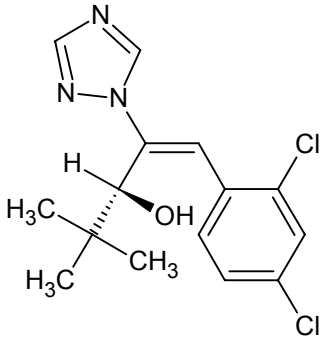
E dimoxystrobin **F** dimoxystrobine (f)	(*E*)-2-(methoxyimino)-*N*-methyl-2-[α-(2,5-xylyloxy)-*o*-tolyl]acetamide (*E*)-2-(méthoxyimino)-*N*-méthyl-2-[2-(2,5-xylyloxy)-méthylphényl]acétamide (α*E*)-2-[(2,5-dimethylphenoxy)methyl]-α-(methoxyimino)-*N*-methylbenzeneacetamide			**F**
		$C_{19}H_{22}N_2O_3$	149961-52-4	
	InChI=1/C19H22N2O3/c1-13-9-10-14(2)17(11-13)24-12-15-7-5-6-8-16(15)18(21-23-4)19(22)20-3/h5-11H,12H2,1-4H3,(H,20,22)/b21-18+/f/h20H			
E diniconazole **F** diniconazole (m)	(*E*)-(*RS*)-1-(2,4-dichlorophenyl)-4,4-dimethyl-2-(1*H*-1,2,4-triazol-1-yl)pent-1-en-3-ol (*E*)-(*RS*)-1-(2,4-dichlorophényl)-4,4-diméthyl-2-(1*H*-1,2,4-triazol-1-yl)=pent-1-én-3-ol (β*E*)-β-[(2,4-dichlorophenyl)methylene]-α-(1,1-dimethylethyl)-1*H*-1,2,4-triazole-1-ethanol			**F**
		$C_{15}H_{17}Cl_2N_3O$	83657-24-3	
	InChI=1/C15H17Cl2N3O/c1-15(2,3)14(21)13(20-9-18-8-19-20)6-10-4-5-11(16)7-12(10)17/h4-9,14,21H,1-3H3/b13-6+/t14-/s3			
E diniconazole-M **F** diniconazole-M (m)	(*E*)-(*R*)-1-(2,4-dichlorophenyl)-4,4-dimethyl-2-(1*H*-1,2,4-triazol-1-yl)pent-1-en-3-ol (*E*)-(*R*)-1-(2,4-dichlorophényl)-4,4-diméthyl-2-(1*H*-1,2,4-triazol-1-yl)=pent-1-én-3-ol (α*R*,β*E*)-β-[(2,4-dichlorophenyl)methylene]-α-(1,1-dimethylethyl)-1*H*-1,2,4-triazole-1-ethanol			**F**
		$C_{15}H_{17}Cl_2N_3O$	83657-18-5	
	InChI=1/C15H17Cl2N3O/c1-15(2,3)14(21)13(20-9-18-8-19-20)6-10-4-5-11(16)7-12(10)17/h4-9,14,21H,1-3H3/b13-6+/t14-/m0/s1			

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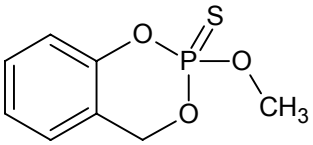
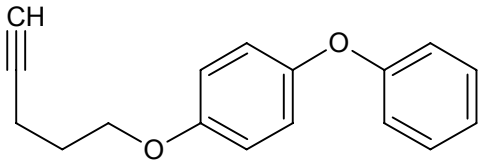
E dioxabenzofos **F** dioxabenzofos (m)	(*RS*)-2-methoxy-4*H*-1,3,2λ^5-benzodioxaphosphinine 2-sulfide			**I**
	2-sulfure de (*RS*)-2-méthoxy-4*H*-1,3,2λ^5,-benzodioxaphosphinine			
	2-methoxy-4*H*-1,3,2-benzodioxaphosphorin 2-sulfide	$C_8H_9O_3PS$	3811-49-2	
	InChI=1/C8H9O3PS/c1-9-12(13)10-6-7-4-2-3-5-8(7)11-12/h2-5H,6H2,1H3/t12-/s3			
E dofenapyn **F** dofénapyne (m)	4-(pent-4-ynyloxy)phenyl phenyl ether			**A**
	éther (ou oxyde) de pent-4-ynyle et de 4-phénoxyphényle			
	1-(4-pentynyloxy)-4-phenoxybenzene	$C_{17}H_{16}O_2$	42873-80-3	
	InChI=1/C17H16O2/c1-2-3-7-14-18-15-10-12-17(13-11-15)19-16-8-5-4-6-9-16/h1,4-6,8-13H,3,7,14H2			

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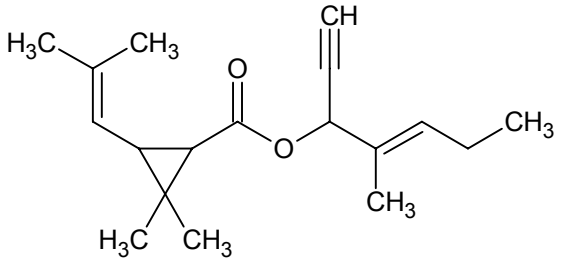
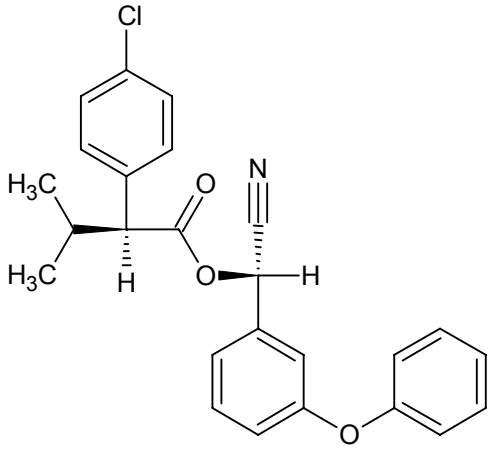
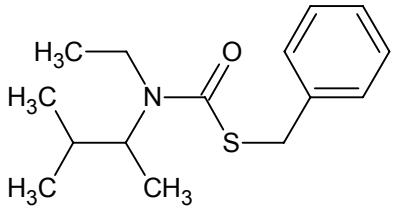
E empenthrin **F** empenthrine (f)	(*E*)-(*RS*)-1-ethynyl-2-methylpent-2-enyl (1*RS*,3*RS*;1*RS*,3*SR*)-2,2-dimethyl-3-(2-methylprop-1-enyl)cyclopropanecarboxylate Alternative Rothamsted-style stereodescriptors: (*E*)-(*RS*)-1-ethynyl-2-methylpent-2-enyl (1*RS*)-*cis-trans*-2,2-dimethyl-3-(2-methylprop-1-enyl)cyclopropanecarboxylate (1*RS*,3*RS*;1*RS*,3*SR*)-2,2-diméthyl-3-(2-méthylprop-1-ènyl)-cyclopropanecarboxylate de (*E*)-(*RS*)-4-méthylhept-1-yn-4-èn-3-yle (2*E*)-1-ethynyl-2-methyl-2-pentenyl 2,2-dimethyl-3-(2-methyl-1-propenyl)cyclopropanecarboxylate			**I**
		$C_{18}H_{26}O_2$	54406-48-3	
	InChI=1/C18H26O2/c1-8-10-13(5)15(9-2)20-17(19)16-14(11-12(3)4)18(16,6)7/h2,10-11,14-16H,8H2,1,3-7H3/b13-10+/t14?,15?,16?			
E esfenvalerate **F** esfenvalérate (m)	(*S*)-α-cyano-3-phenoxybenzyl (*S*)-2-(4-chlorophenyl)-3-methylbutyrate (*S*)-2-(4-chlorophényl)isovalérate de (*S*)-cyano-(3-phénoxyphényl)méthyle (*S*)-cyano(3-phenoxyphenyl)methyl (α*S*)-4-chloro-α-(1-methylethyl)benzeneacetate			**I**
		$C_{25}H_{22}ClNO_3$	66230-04-4	
	InChI=1/C25H22ClNO3/c1-17(2)24(18-11-13-20(26)14-12-18)25(28)30-23(16-27)19-7-6-10-22(15-19)29-21-8-4-3-5-9-21/h3-15,17,23-24H,1-2H3/t23-,24+/m1/s1			
E esprocarb **F** esprocarb (m)	*S*-benzyl (*RS*)-1,2-dimethylpropyl(ethyl)thiocarbamate (*RS*)-éthyl-(3-méthylbut-2-yl)thiocarbamate de *S*-benzyle *S*-(phenylmethyl) (1,2-dimethylpropyl)ethylcarbamothioate			**H**
		$C_{15}H_{23}NOS$	85785-20-2	
	InChI=1/C15H23NOS/c1-5-16(13(4)12(2)3)15(17)18-11-14-9-7-6-8-10-14/h6-10,12-13H,5,11H2,1-4H3/t13-/s3			

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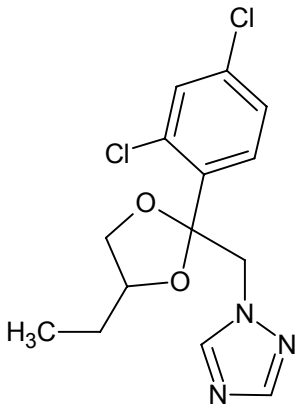
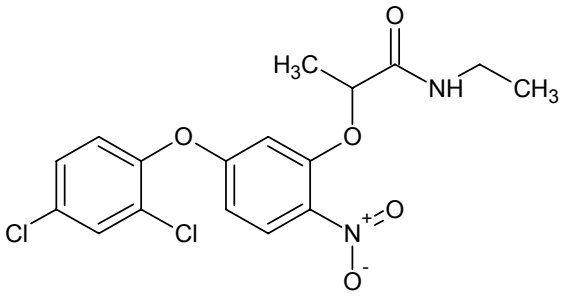
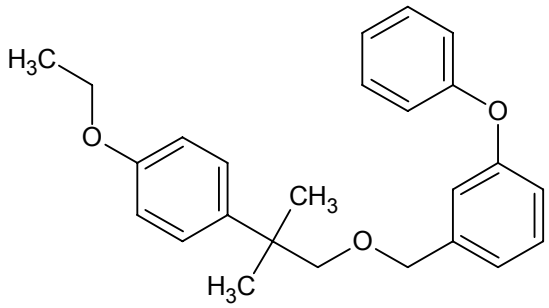
E etaconazole **F** étaconazole (m)	1-[(2*RS*,4*RS*;2*RS*,4*SR*)-2-(2,4-dichlorophenyl)-4-ethyl-1,3-dioxolan-2-ylmethyl]-1*H*-1,2,4-triazole 1-[(2*RS*,4*RS*;2*RS*,4*SR*)-2-(2,4-dichlorophényl)-4-éthyl-1,3-dioxolan-2-ylméthyl]-1*H*-1,2,4-triazole 1-[[2-(2,4-dichlorophenyl)-4-ethyl-1,3-dioxolan-2-yl]methyl]-1*H*-1,2,4-triazole	 $C_{14}H_{15}Cl_2N_3O_2$	60207-93-4	**F**
	InChI=1/C14H15Cl2N3O2/c1-2-11-6-20-14(21-11,7-19-9-17-8-18-19)12-4-3-10(15)5-13(12)16/h3-5,8-9,11H,2,6-7H2,1H3/t11?,14?			
	NOTE The ratio of isomers should be stated. NOTE Il convient de préciser la proportion des isomères.			
E etnipromid **F** etnipromide (m)	(*RS*)-2-[5-(2,4-dichlorophenoxy)-2-nitrophenoxy]-*N*-ethylpropionamide (*RS*)-2-[5-(2,4-dichlorophénoxy)-2-nitrophénoxy]-*N*-éthylpropionamide 2-[5-(2,4-dichlorophenoxy)-2-nitrophenoxy]-*N*-ethylpropanamide	 $C_{17}H_{16}Cl_2N_2O_5$	76120-02-0	**H**
	InChI=1/C17H16Cl2N2O5/c1-3-20-17(22)10(2)25-16-9-12(5-6-14(16)21(23)24)26-15-7-4-11(18)8-13(15)19/h4-10H,3H2,1-2H3,(H,20,22)/t10-/s3/f/h20H			
E etofenprox **F** étofénoprox (m)	2-(4-ethoxyphenyl)-2-methylpropyl 3-phenoxybenzyl ether éther (ou oxyde) de 2-(4-éthoxyphényl)-2-méthylpropyle et de 3-phénoxyphénylméthyle 1-[[2-(4-ethoxyphenyl)-2-methylpropoxy]methyl]-3-phenoxybenzene	 $C_{25}H_{28}O_3$	80844-07-1	**I**
	InChI=1/C25H28O3/c1-4-27-22-15-13-21(14-16-22)25(2,3)19-26-18-20-9-8-12-24(17-20)28-23-10-6-5-7-11-23/h5-17H,4,18-19H2,1-3H3			

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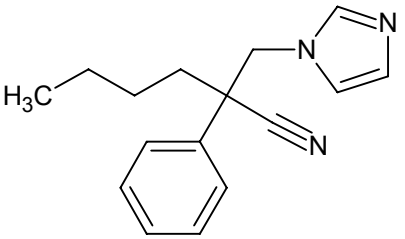
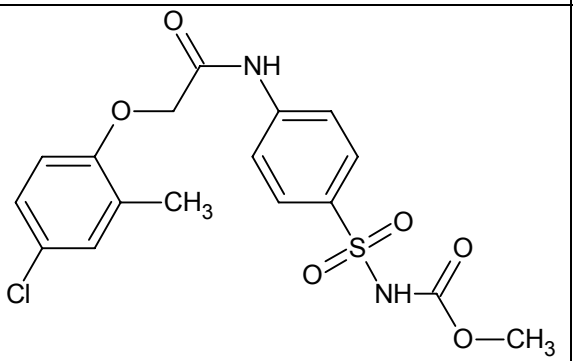
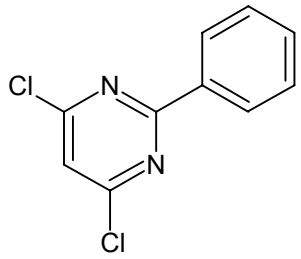
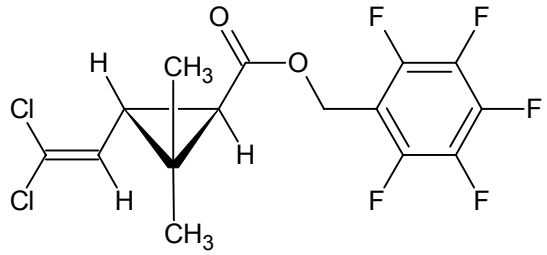
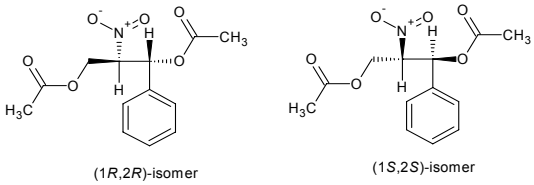
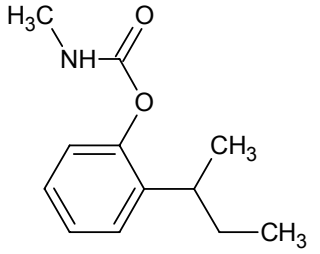
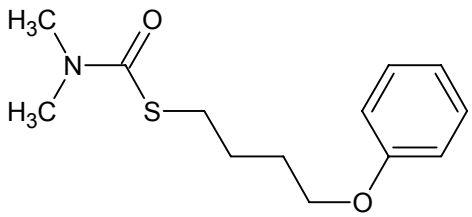
	IUPAC International Chemical Identifier (InChI™)			
E fenapanil **F** fénapanil (m)	(*RS*)-2-(imidazol-1-ylmethyl)-2-phenylhexanenitrile (*RS*)-2-(imidazol-1-ylméthyl)-2-phénylhexanenitrile α-butyl-α-phenyl-1*H*-imidazole-1-propanenitrile			**F**
		$C_{16}H_{19}N_3$	61019-78-1	
	InChI=1/C16H19N3/c1-2-3-9-16(12-17,13-19-11-10-18-14-19)15-7-5-4-6-8-15/h4-8,10-11,14H,2-3,9,13H2,1H3/t16-/s3			
E fenasulam **F** fénasulame (m)	methyl 4-[2-(4-chloro-*o*-tolyloxy)=acetamido]phenylsulfonylcarbamate Alternatively: methyl (*N*-4-chloro-*o*-tolyloxyacetylsulfanilyl)carbamate 4-[(4-chloro-2-méthylphénoxy)=acétamido]phénylsulfonylcarbamate de méthyle methyl [[4-[[(4-chloro-2-methylphenoxy)acetyl]amino]phenyl]=sulfonyl]carbamate			**H**
		$C_{17}H_{17}ClN_2O_6S$	78357-48-9	
	InChI=1/C17H17ClN2O6S/c1-11-9-12(18)3-8-15(11)26-10-16(21)19-13-4-6-14(7-5-13)27(23,24)20-17(22)25-2/h3-9H,10H2,1-2H3,(H,19,21)(H,20,22)/f/h19-20H			
E fenclorim **F** fenclorime (f)	4,6-dichloro-2-phenylpyrimidine 4,6-dichloro-2-phénylpyrimidine 4,6-dichloro-2-phenylpyrimidine			**S**
		$C_{10}H_6Cl_2N_2$	3740-92-9	
	InChI=1/C10H6Cl2N2/c11-8-6-9(12)14-10(13-8)7-4-2-1-3-5-7/h1-6H			

E: Common name F: Nom commun	Chemical name Nom chimique E: IUPAC F: UICPA C: CAS	Structure Structure Molecular formula Formule brute	CAS Registry Number® Numéro d'enregistrement CAS®	Use Application
	IUPAC International Chemical Identifier (InChI™)			
E fenfluthrin **F** fenfluthrine (f)	2,3,4,5,6-pentafluorobenzyl (1*R*,3*S*)-3-(2,2-dichlorovinyl)-2,2-dimethylcyclopropanecarboxylate Alternative Rothamsted-style stereodescriptors: 2,3,4,5,6-pentafluorobenzyl (1*R*)-*trans*-3-(2,2-dichlorovinyl)-2,2-dimethylcyclopropanecarboxylate (1*R*,3*S*)-3-(2,2-dichlorovinyl)-2,2-diméthylcyclopropanecarboxylate de pentafluorobenzyle (pentafluorophenyl)methyl (1*R*,3*S*)-3-(2,2-dichloroethenyl)-2,2-dimethylcyclopropanecarboxylate	 $C_{15}H_{11}Cl_2F_5O_2$	75867-00-4	I
	InChI=1/C15H11Cl2F5O2/c1-15(2)6(3-7(16)17)8(15)14(23)24-4-5-9(18)11(20)13(22)12(21)10(5)19/h3,6,8H,4H2,1-2H3/t6-,8+/m1/s1			
E fenitropan **F** fénitropane (m)	(1*RS*,2*RS*)-2-nitro-1-phenyltrimethylene di(acetate) diacétate de (1*RS*,2*RS*)-2-nitro-1-phénylprop-1,3-diyle (1*R*,2*R*)-*rel*-2-nitro-1-phenyl-1,3-propanediyl diacetate	 (1*R*,2*R*)-isomer (1*S*,2*S*)-isomer $C_{13}H_{15}NO_6$	65934-95-4	F
	InChI=1/C13H15NO6/c1-9(15)19-8-12(14(17)18)13(20-10(2)16)11-6-4-3-5-7-11/h3-7,12-13H,8H2,1-2H3/t12-,13-/s3			
E fenobucarb **F** fénobucarb (m)	(*RS*)-2-*sec*-butylphenyl methylcarbamate méthylcarbamate de (*RS*)-2-(but-2-yl)-phényle 2-(1-methylpropyl)phenyl methylcarbamate	 $C_{12}H_{17}NO_2$	3766-81-2	I
	InChI=1/C12H17NO2/c1-4-9(2)10-7-5-6-8-11(10)15-12(14)13-3/h5-9H,4H2,1-3H3,(H,13,14)/t9-/s3/f/h13H			
E fenothiocarb **F** fénothiocarb (m)	*S*-4-phenoxybutyl dimethyl(thiocarbamate) diméthylthiocarbamate de *S*-4-phénoxybutyle *S*-(4-phenoxybutyl) dimethylcarbamothioate	 $C_{13}H_{19}NO_2S$	62850-32-2	A
	InChI=1/C13H19NO2S/c1-14(2)13(15)17-11-7-6-10-16-12-8-4-3-5-9-12/h3-5,8-9H,6-7,10-11H2,1-2H3			

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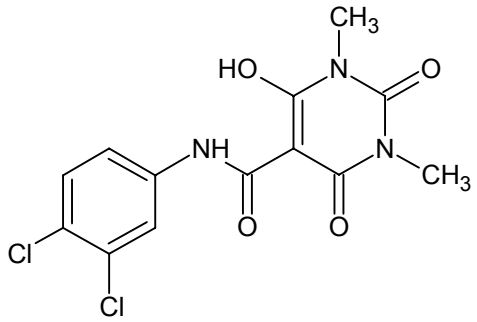
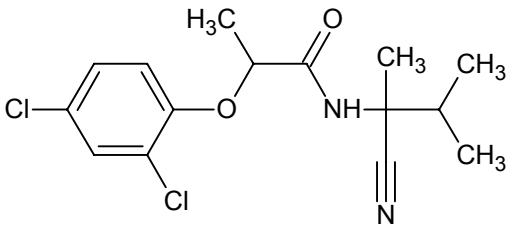
E fenoxacrim **F** fénoxacrim (m)	3′,4′-dichloro-1,2,3,4-tetrahydro-6-hydroxy-1,3-dimethyl-2,4-dioxopyrimidine-5-carboxanilide 3′,4′-dichloro-6-hydroxy-1,3-diméthyl-2,4-dioxo-(1,2,3,4-tétrahydropyrimidine)-5-carboxanilide *N*-(3,4-dichlorophenyl)hexahydro-1,3-dimethyl-2,4,6-trioxo-5-pyrimidinecarboxamide			**I**
		$C_{13}H_{11}Cl_2N_3O_4$	65400-98-8	
	InChI=1/C13H11Cl2N3O4/c1-17-11(20)9(12(21)18(2)13(17)22)10(19)16-6-3-4-7(14)8(15)5-6/h3-5,20H,1-2H3,(H,16,19)/f/h16H CAS-preferred tautomer InChI=1/C13H11Cl2N3O4/c1-17-11(20)9(12(21)18(2)13(17)22)10(19)16-6-3-4-7(14)8(15)5-6/h3-5,9H,1-2H3,(H,16,19)/f/h16H			
	NOTE The CAS name and Registry Number are for the tautomer that is preferred under the CAS nomenclature rules. NOTE Le nom et le numéro du CAS font référence au tautomère qui est préféré selon les règles de nomenclature du CAS.			
E fenoxanil **F** fénoxanile (m)	mixture of 85 % (*R*)-*N*-[(*RS*)-1-cyano-1,2-dimethylpropyl]-2-(2,4-dichlorophenoxy)propionamide and 15 % (*S*)-*N*-[(*RS*)-1-cyano-1,2-dimethylpropyl]-2-(2,4-dichlorophenoxy)propionamide mélange de 85 % de (*R*)-*N*-[(*RS*)-2-cyano-3-méthylbut-2-yl]-2-(2,4-dichlorophénoxy)propionamide et de 15% de (*S*)-*N*-[(*RS*)-2-cyano-3-méthylbut-2-yl]-2-(2,4-dichlorophénoxy)propionamide *N*-(1-cyano-1,2-dimethylpropyl)-2-(2,4-dichlorophenoxy)propanamide			**F**
		$C_{15}H_{18}Cl_2N_2O_2$	115852-48-7	
	InChI=1/C15H18Cl2N2O2/c1-9(2)15(4,8-18)19-14(20)10(3)21-13-6-5-11(16)7-12(13)17/h5-7,9-10H,1-4H3,(H,19,20)/t10?,15?/f/h19H			

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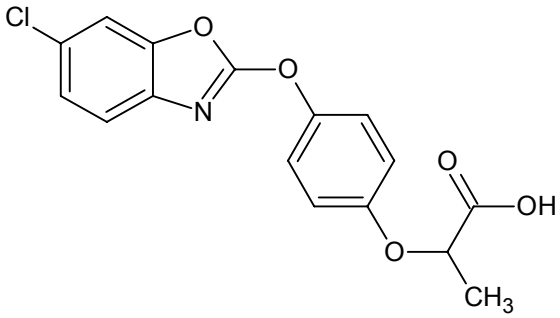
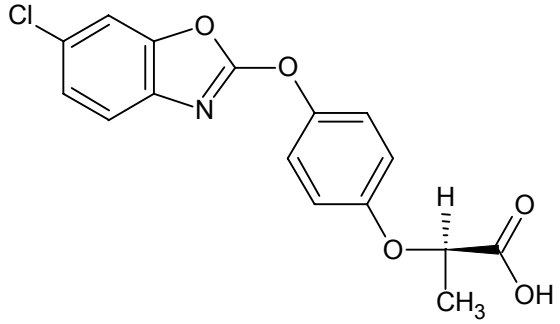
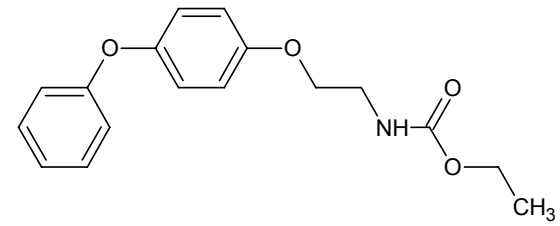
E fenoxaprop **F** fénoxaprop (m)	(*RS*)-2-[4-(6-chloro-1,3-benzoxazol-2-yloxy)phenoxy]propionic acid acide (*RS*)-2-[4-(6-chloro-1,3-benzoxazol-2-yloxy)phénoxy]propionique 2-[4-[(6-chloro-2-benzoxazolyl)oxy]phenoxy]propanoic acid			**H**
		$C_{16}H_{12}ClNO_5$	95617-09-7	
	InChI=1/C16H12ClNO5/c1-9(15(19)20)21-11-3-5-12(6-4-11)22-16-18-13-7-2-10(17)8-14(13)23-16/h2-9H,1H3,(H,19,20)/t9-/s3/f/h19H fenoxaprop-ethyl InChI=1/C18H16ClNO5/c1-3-22-17(21)11(2)23-13-5-7-14(8-6-13)24-18-20-15-9-4-12(19)10-16(15)25-18/h4-11H,3H2,1-2H3/t11-/s3			
	NOTE It should be stated which ester or salt is present, for example fenoxaprop-ethyl [66441-23-4]. NOTE Il convient de préciser quel est l'ester ou le sel présent, par exemple fénoxaprop-éthyle [66441-23-4].			
E fenoxaprop-P **F** fénoxaprop-P (m)	(*R*)-2-[4-(6-chloro-1,3-benzoxazol-2-yloxy)phenoxy]propionic acid acide (*R*)-2-[4-(6-chloro-1,3-benzoxazol-2-yloxy)phénoxy]propionique (2*R*)-2-[4-[(6-chloro-2-benzoxazolyl)oxy]phenoxy]propanoic acid			**H**
		$C_{16}H_{12}ClNO_5$	113158-40-0	
	InChI=1/C16H12ClNO5/c1-9(15(19)20)21-11-3-5-12(6-4-11)22-16-18-13-7-2-10(17)8-14(13)23-16/h2-9H,1H3,(H,19,20)/t9-/m1/s1/f/h19H fenoxaprop-P-ethyl InChI=1/C18H16ClNO5/c1-3-22-17(21)11(2)23-13-5-7-14(8-6-13)24-18-20-15-9-4-12(19)10-16(15)25-18/h4-11H,3H2,1-2H3/t11-/m1/s1			
	NOTE It should be stated which ester or salt is present, for example fenoxaprop-P-ethyl [71283-80-2]. NOTE Il convient de préciser quel est l'ester ou le sel présent, par exemple fénoxaprop-P-éthyle [71283-80-2].			
E fenoxycarb **F** fénoxycarb (m)	ethyl 2-(4-phenoxyphenoxy)ethylcarbamate 2-(4-phénoxyphénoxy)éthylcarbamate d'éthyle ethyl [2-(4-phenoxyphenoxy)ethyl]carbamate			**IGR**
		$C_{17}H_{19}NO_4$	79127-80-3	
	InChI=1/C17H19NO4/c1-2-20-17(19)18-12-13-21-14-8-10-16(11-9-14)22-15-6-4-3-5-7-15/h3-11H,2,12-13H2,1H3,(H,18,19)/f/h18H			

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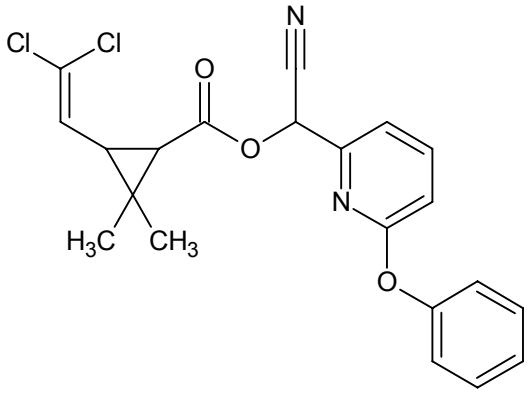
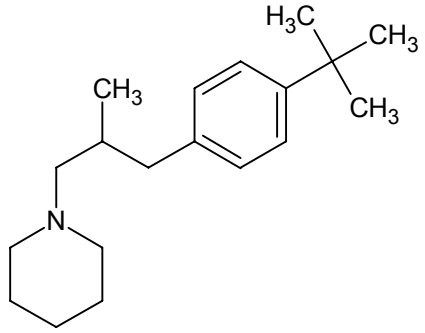
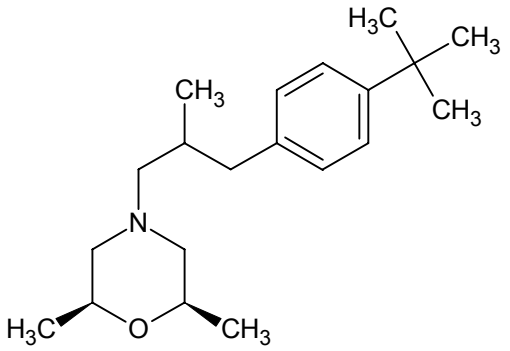
	IUPAC International Chemical Identifier (InChI™)			
E fenpirithrin **F** fenpirithrine (f)	(*RS*)-cyano(6-phenoxy-2-pyridyl)methyl (1*RS*,3*RS*;1*RS*,3*SR*)-3-(2,2-dichlorovinyl)-2,2-dimethylcyclopropanecarboxylate Alternative Rothamsted-style stereodescriptors: (*RS*)-cyano(6-phenoxy-2-pyridyl)methyl (1*RS*)-*cis-trans*-3-(2,2-dichlorovinyl)-2,2-dimethylcyclopropanecarboxylate (*RS*)-(3-(2,2-dichlorovinyl)-2,2-diméthyl)cyclopropanecarboxylate de cyano-(6-phénoxypyrid-2-yl)-méthyle cyano(6-phenoxy-2-pyridinyl)methyl 3-(2,2-dichloroethenyl)-2,2-dimethylcyclopropanecarboxylate		$C_{21}H_{18}Cl_2N_2O_3$	68523-18-2	I
	InChI=1/C21H18Cl2N2O3/c1-21(2)14(11-17(22)23)19(21)20(26)28-16(12-24)15-9-6-10-18(25-15)27-13-7-4-3-5-8-13/h3-11,14,16,19H,1-2H3/t14?,16?,19?			
	NOTE The (1*R*,3*R*) to (1*R*,3*S*) and the (1*S*,3*S*) to (1*S*,3*R*) ratios each vary from 40 to 60 and 60 to 40, respectively. NOTE Les proportions entre les racémates (1*R*,3*R*) et (1*R*,3*S*) et entre les racémates (1*S*,3*S*) et (1*S*,3*R*) varient de 40 à 60 et de 60 à 40, respectivement.			
E fenpropidin **F** fenpropidine (f)	(*RS*)-1-[3-(4-*tert*-butylphenyl)-2-methylpropyl]piperidine (*RS*)-*N*-[3-(4-*tert*-butylphényl)-2-méthylpropyl]pipéridine ou (*RS*)-1-[3-(4-*tert*-Butylphényl)-2-méthylpropyl]pipéridine 1-[3-[4-(1,1-dimethylethyl)phenyl]-2-methylpropyl]piperidine		$C_{19}H_{31}N$	67306-00-7	F
	InChI=1/C19H31N/c1-16(15-20-12-6-5-7-13-20)14-17-8-10-18(11-9-17)19(2,3)4/h8-11,16H,5-7,12-15H2,1-4H3/t16-/s3			
E fenpropimorph **F** fenpropimorphe (m)	*cis*-4-[(*RS*)-3-(4-*tert*-butylphenyl)-2-methylpropyl]-2,6-dimethylmorpholine *cis*-4-[(*RS*)-3-(4-*tert*-butylphényl)-2-méthylpropyl]-2,6-diméthylmorpholine ou *cis*-*N*-[(*RS*)-3-(4-*tert*-butylphényl)-2-méthylpropyl]-2,6-diméthylmorpholine (2*R*,6*S*)-*rel*-4-[3-[4-(1,1-dimethylethyl)phenyl]-2-methylpropyl]-2,6-dimethylmorpholine		$C_{20}H_{33}NO$	67564-91-4	F
	InChI=1/C20H33NO/c1-15(12-21-13-16(2)22-17(3)14-21)11-18-7-9-19(10-8-18)20(4,5)6/h7-10,15-17H,11-14H2,1-6H3/t15?,16-,17+			

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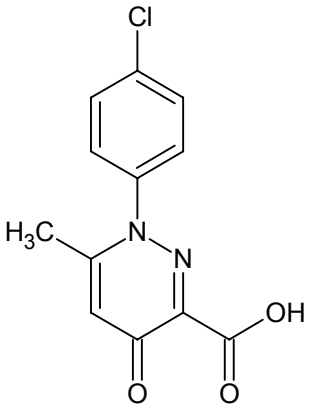
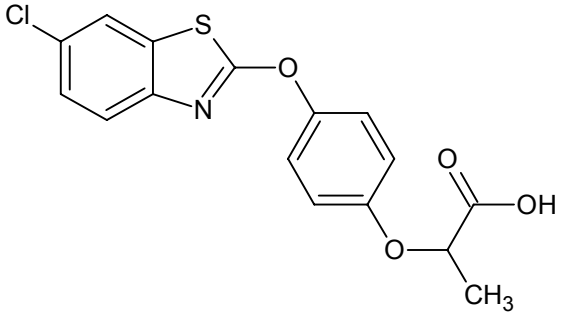
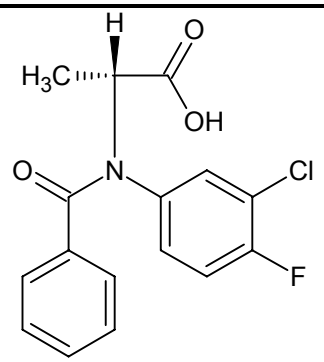
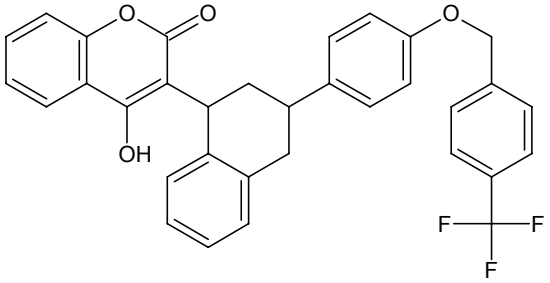
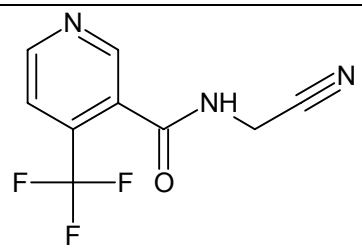
		Molecular formula Formule brute	CAS Registry Number® Numéro d'enregistrement CAS®	Applic-ation
	IUPAC International Chemical Identifier (InChI™)			
E fenridazon **F** fenridazon (m)	1-(4-chlorophenyl)-1,4-dihydro-6-methyl-4-oxopyridazine-3-carboxylic acid acide 1-(4-chlorophényl)-6-méthyl-4-oxo-1,4-dihydropyridazine-3-carboxylique 1-(4-chlorophenyl)-1,4-dihydro-6-methyl-4-oxo-3-pyridazinecarboxylic acid	 $C_{12}H_9ClN_2O_3$	68254-10-4	**P**
	InChI=1/C12H9ClN2O3/c1-7-6-10(16)11(12(17)18)14-15(7)9-4-2-8(13)3-5-9/h2-6H,1H3,(H,17,18)/f/h17H fenridazon-potassium InChI=1/C12H9ClN2O3.K/c1-7-6-10(16)11(12(17)18)14-15(7)9-4-2-8(13)3-5-9;/h2-6H,1H3,(H,17,18);/q;+1/p-1/fC12H8ClN2O3.K/q-1;m			
	NOTE 1 It should be stated which ester or salt is present, for example fenridazon-potassium [83588-43-6]. NOTE 1 Il convient de préciser quel est l'ester ou le sel présent, par exemple fenridazon-potassium [83588-43-6]. NOTE 2 The name “fenridazon” is not acceptable for use in Hungary because of the risk of confusion with the trade name “Rideon”. NOTE 2 Le nom «fenridazon» n'est pas acceptable pour l'emploi en Hongrie, car il entre en conflit avec le nom commercial «Rideon».			
E fenthiaprop **F** fentiaprop (m)	(*RS*)-2-[4-(6-chloro-1,3-benzothiazol-2-yloxy)phenoxy]propionic acid acide (*RS*)-2-[4-(6-chloro-1,3-benzothiazol-2-yloxy)phénoxy]propionique 2-[4-[(6-chloro-2-benzothiazolyl)oxy]=phenoxy]propanoic acid	 $C_{16}H_{12}ClNO_4S$	73519-50-3	**H**
	InChI=1/C16H12ClNO4S/c1-9(15(19)20)21-11-3-5-12(6-4-11)22-16-18-13-7-2-10(17)8-14(13)23-16/h2-9H,1H3,(H,19,20)/t9-/s3/f/h19H fenthiaprop-ethyl InChI=1/C18H16ClNO4S/c1-3-22-17(21)11(2)23-13-5-7-14(8-6-13)24-18-20-15-9-4-12(19)10-16(15)25-18/h4-11H,3H2,1-2H3/t11-/s3			
	NOTE It should be stated which ester or salt is present, for example fenthiaprop-ethyl [66441-11-0]. NOTE Il convient de préciser quel est l'ester ou le sel présent, par exemple fentiaprop-éthyle [66441-11-0].			

E: Common name F: Nom commun	Chemical name Nom chimique E: IUPAC F: UICPA C: CAS	Structure Structure Molecular formula Formule brute	CAS Registry Number® Numéro d'enregistrement CAS®	Use Applic-ation
	IUPAC International Chemical Identifier (InChI™)			
E flamprop-M **F** flamprop-M (m)	*N*-benzoyl-*N*-(3-chloro-4-fluorophenyl)-D-alanine *N*-(3-chloro-4-fluorophényl)-*N*-benzoyl-D-alanine *N*-benzoyl-*N*-(3-chloro-4-fluorophenyl)-D-alanine			**H**
		$C_{16}H_{13}ClFNO_3$	90134-59-1	
	InChI=1/C16H13ClFNO3/c1-10(16(21)22)19(12-7-8-14(18)13(17)9-12)15(20)11-5-3-2-4-6-11/h2-10H,1H3,(H,21,22)/t10-/m1/s1/f/h21H flamprop-M-isopropyl InChI=1/C19H19ClFNO3/c1-12(2)25-19(24)13(3)22(15-9-10-17(21)16(20)11-15)18(23)14-7-5-4-6-8-14/h4-13H,1-3H3/t13-/m1/s1 flamprop-M-methyl InChI=1/C17H15ClFNO3/c1-11(17(22)23-2)20(13-8-9-15(19)14(18)10-13)16(21)12-6-4-3-5-7-12/h3-11H,1-2H3/t11-/m1/s1			
	NOTE It should be stated which ester or salt is present, for example flamprop-M-isopropyl [63782-90-1] or flamprop-M-methyl [63729-98-6]. NOTE Il convient de préciser quel est l'ester ou le sel présent, par exemple flamprop-M-isopropyle [63782-90-1] ou flamprop-M-méthyle [63729-98-6].			
E flocoumafen **F** flocoumafène (m)	4-hydroxy-3-[(1*RS*,3*RS*;1*RS*,3*SR*)-1,2,3,4-tetrahydro-3-[4-(4-trifluoromethylbenzyloxy)phenyl]-1-naphthyl]coumarin (mixture of *cis*- to *trans*- isomers in the ratio range 60:40 to 40:60 respectively) mélange des isomères (1*R*,3*R*) et (1*R*,3*S*) de la 4-hydroxy-3-[1,2,3,4-tétrahydro-3-((4-trifluorométhylbenzyloxy)phényl]-napht-1-ylcoumarine en proportions allant de 60/40 à 40/60 4-hydroxy-3-[1,2,3,4-tetrahydro-3-[4-[[4-(trifluoromethyl)phenyl]methoxy]=phenyl]-1-naphthalenyl]-2*H*-1-benzopyran-2-one			**R**
		$C_{33}H_{25}F_3O_4$	90035-08-8	
	InChI=1/C33H25F3O4/c34-33(35,36)24-13-9-20(10-14-24)19-39-25-15-11-21(12-16-25)23-17-22-5-1-2-6-26(22)28(18-23)30-31(37)27-7-3-4-8-29(27)40-32(30)38/h1-16,23,28,37H,17-19H2/t23?,28?			
E flonicamid **F** flonicamide (m)	*N*-cyanomethyl-4-(trifluoromethyl)nicotinamide *N*-cyanométhyl-4-(trifluorométhyl)nicotinamide *N*-(cyanomethyl)-4-(trifluoromethyl)-3-pyridinecarboxamide			**I**
		$C_9H_6F_3N_3O$	158062-67-0	
	InChI=1/C9H6F3N3O/c10-9(11,12)7-1-3-14-5-6(7)8(16)15-4-2-13/h1,3,5H,4H2,(H,15,16)/f/h15H			

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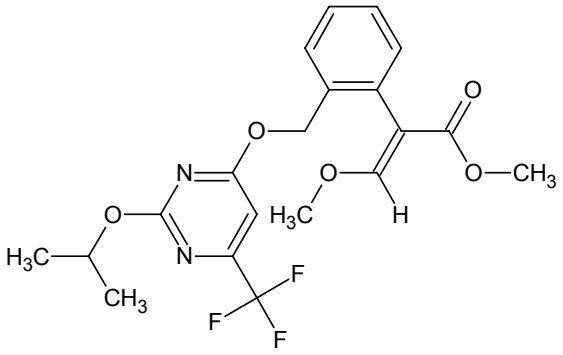
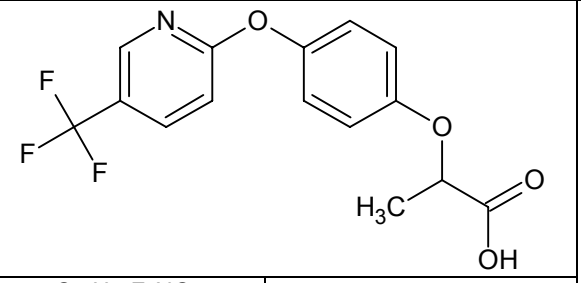
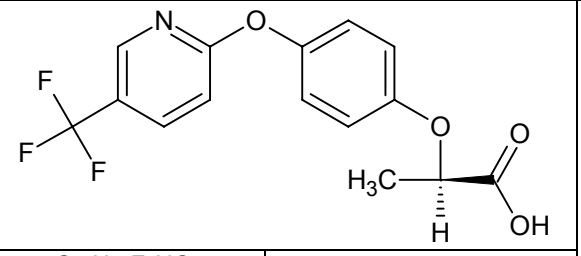
E fluacrypyrim **F** fluacrypyrime (m)	methyl (*E*)-2-{α-[2-isopropoxy-6-(trifluoromethyl)pyrimidin-4-yloxy]-*o*-tolyl}-3-methoxyacrylate (*E*)-2-{2-[2-isopropoxy-6-(trifluorométhyl)pyrimidin-4-yloxy]-méthyl}phényl-3-méthoxyacrylate de méthyle methyl (α*E*)-α-(methoxymethylene)-2-[[[2-(1-methylethoxy)-6-trifluoromethyl)-4-pyrimidinyl]oxy]methyl]benzeneacetate			**A**
		$C_{20}H_{21}F_3N_2O_5$	229977-93-9	
	InChI=1/C20H21F3N2O5/c1-12(2)30-19-24-16(20(21,22)23)9-17(25-19)29-10-13-7-5-6-8-14(13)15(11-27-3)18(26)28-4/h5-9,11-12H,10H2,1-4H3/b15-11+			
E fluazifop **F** fluazifop (m)	(*RS*)-2-{4-[5-(trifluoromethyl)-2-pyridyloxy]phenoxy}propionic acid acide (*RS*)-2-{4-[5-(trifluorométhyl)pyrid-2-yloxy]phénoxy}propionique 2-[4-[[5-(trifluoromethyl)-2-pyridinyl]oxy]phenoxy]propanoic acid			**H**
		$C_{15}H_{12}F_3NO_4$	69335-91-7	
	InChI=1/C15H12F3NO4/c1-9(14(20)21)22-11-3-5-12(6-4-11)23-13-7-2-10(8-19-13)15(16,17)18/h2-9H,1H3,(H,20,21)/t9-/s3/f/h20H fluazifop-butyl InChI=1/C19H20F3NO4/c1-3-4-11-25-18(24)13(2)26-15-6-8-16(9-7-15)27-17-10-5-14(12-23-17)19(20,21)22/h5-10,12-13H,3-4,11H2,1-2H3/t13-/s3			
	NOTE It should be stated which ester or salt is present, for example fluazifop-butyl [69806-50-4]. NOTE Il convient de préciser quel est l'ester ou le sel présent, par exemple fluazifop-butyl [69806-50-4].			
E fluazifop-P **F** fluazifop-P (m)	(*R*)-2-{4-[5-(trifluoromethyl)-2-pyridyloxy]phenoxy}propionic acid acide (*R*)-2-{4-[5-(trifluorométhyl)pyrid-2-yloxy]phénoxy}propionique (2*R*)-2-[4-[[5-(trifluoromethyl)-2-pyridinyl]oxy]phenoxy]propanoic acid			**H**
		$C_{15}H_{12}F_3NO_4$	83066-88-0	
	InChI=1/C15H12F3NO4/c1-9(14(20)21)22-11-3-5-12(6-4-11)23-13-7-2-10(8-19-13)15(16,17)18/h2-9H,1H3,(H,20,21)/t9-/m1/s1/f/h20H fluazifop-P-butyl InChI=1/C19H20F3NO4/c1-3-4-11-25-18(24)13(2)26-15-6-8-16(9-7-15)27-17-10-5-14(12-23-17)19(20,21)22/h5-10,12-13H,3-4,11H2,1-2H3/t13-/m1/s1			
	NOTE It should be stated which ester or salt is present, for example fluazifop-P-butyl [79241-46-6]. NOTE Il convient de préciser quel est l'ester ou le sel présent, par exemple fluazifop-P-butyl [79241-46-6].			

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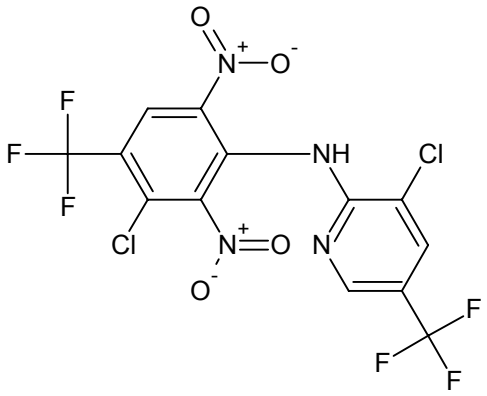
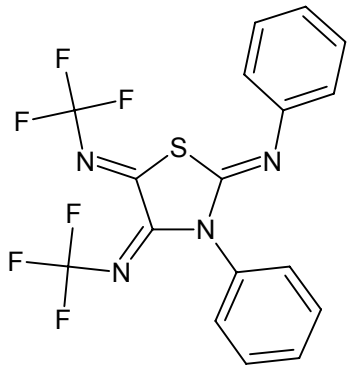
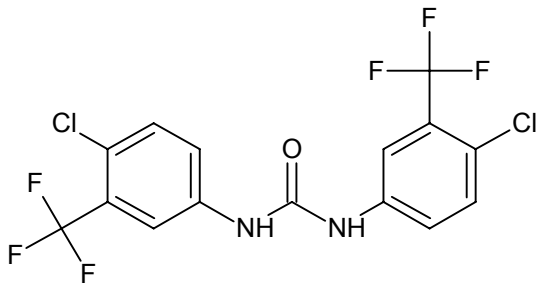
E fluazinam **F** fluaziname (m)	3-chloro-*N*-(3-chloro-5-trifluoromethyl-2-pyridyl)-α,α,α-trifluoro-2,6-dinitro-*p*-toluidine 3-chloro-*N*-(3-chloro-5-trifluorométhylpyrid-2-yl)-4-trifluorométhyl-2,6-dinitroaniline 3-chloro-*N*-[3-chloro-2,6-dinitro-4-(trifluoromethyl)phenyl]-5-(trifluoromethyl)-2-pyridinamine			**F**
		$C_{13}H_4Cl_2F_6N_4O_4$	79622-59-6	
	InChI=1/C13H4Cl2F6N4O4/c14-6-1-4(12(16,17)18)3-22-11(6)23-9-7(24(26)27)2-5(13(19,20)21)8(15)10(9)25(28)29/h1-3H,(H,22,23)/f/h23H			
E flubenzimine **F** flubenzimine (f)	(2*Z*,4*E*,5*Z*)-N^2,3-diphenyl-N^4,N^5-bis(trifluoromethyl)-1,3-thiazolidine-2,4,5-triimine (2*Z*,4*E*,5*Z*)-3-phényl-2-phénylimino-4,5-bis((trifluorométhyl)imino)-thiazolidine [N(*Z*)]-*N*-[(4*E*,5*Z*)-3-phenyl-4,5-bis[(trifluoromethyl)imino]-2-thiazolidinylidene]benzenamine			**A**
		$C_{17}H_{10}F_6N_4S$	37893-02-0	
	InChI=1/C17H10F6N4S/c18-16(19,20)25-13-14(26-17(21,22)23)28-15(24-11-7-3-1-4-8-11)27(13)12-9-5-2-6-10-12/h1-10H/b24-15-,25-13+,26-14-			
E flucofuron **F** flucofuron (m)	1,3-bis(4-chloro-α,α,α-trifluoro-*m*-tolyl)urea 1,3-bis(4-chloro-3-trifluorométhylphényl)urée *N,N'*-bis[4-chloro-3-(trifluoromethyl)phenyl]urea			**I**
		$C_{15}H_8Cl_2F_6N_2O$	370-50-3	
	InChI=1/C15H8Cl2F6N2O/c16-11-3-1-7(5-9(11)14(18,19)20)24-13(26)25-8-2-4-12(17)10(6-8)15(21,22)23/h1-6H,(H2,24,25,26)/f/h24-25H			

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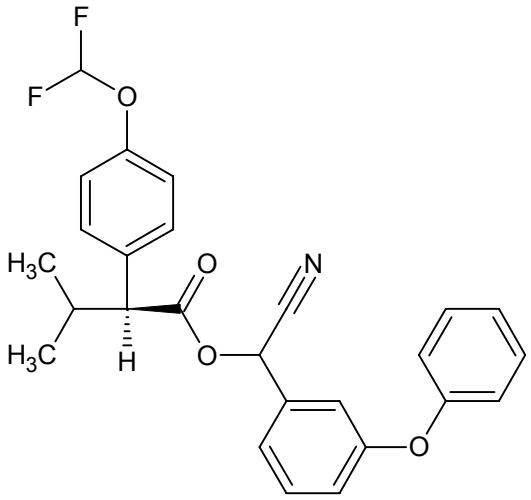
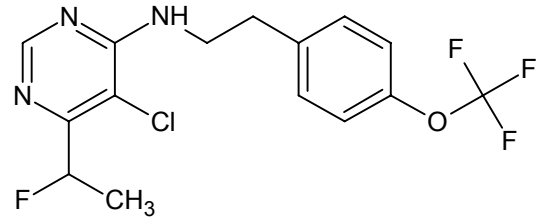
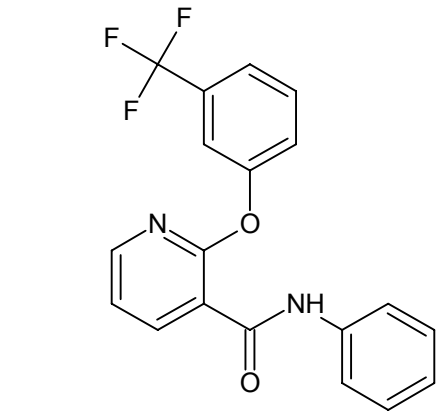
		Molecular formula Formule brute	CAS Registry Number® Numéro d'enregistrement CAS®	Applic-ation
	IUPAC International Chemical Identifier (InChI™)			
E flucythrinate **F** flucythrinate (m)	(*RS*)-α-cyano-3-phenoxybenzyl (*S*)-2-(4-difluoromethoxyphenyl)-3-methylbutyrate (*RS*)-2(4-difluorométhoxyphényl)-3-méthylbutyrate de cyano(3-phénoxyphényl)méthyle cyano(3-phenoxyphenyl)methyl (α*S*)-4-(difluoromethoxy)-α-(1-methylethyl)benzeneacetate			**A** **I**
		$C_{26}H_{23}F_2NO_4$	70124-77-5	
	InChI=1/C26H23F2NO4/c1-17(2)24(18-11-13-21(14-12-18)32-26(27)28)25(30)33-23(16-29)19-7-6-10-22(15-19)31-20-8-4-3-5-9-20/h3-15,17,23-24,26H,1-2H3/t23?,24-/m0/s1			
E flufenerim **F** flufénérime (f)	{5-chloro-6-[(*RS*)-1-fluoroethyl]pyrimidin-4-yl}[4-(trifluoromethoxy)phenethyl]amine (*RS*)-*N*-[5-chloro-6-(1-fluoroéthyl)pyrimidin-4-yl]-4-trifluorométhoxy)phénéthylamine 5-chloro-6-(1-fluoroethyl)-*N*-[2-[4-(trifluoromethoxy)phenyl]ethyl]-4-pyrimidinamine			**I**
		$C_{15}H_{14}ClF_4N_3O$	170015-32-4	
	InChI=1/C15H14ClF4N3O/c1-9(17)13-12(16)14(23-8-22-13)21-7-6-10-2-4-11(5-3-10)24-15(18,19)20/h2-5,8-9H,6-7H2,1H3,(H,21,22,23)/t9-/s3/f/h21H			
E flufenican **F** flufénican (m)	2-(α,α,α-trifluoro-*m*-tolyloxy)nicotinanilide 2-(3-trifluorométhylphénoxy)nicotinanilide *N*-phenyl-2-[3-(trifluoromethyl)phenoxy]-3-pyridinecarboxamide			**H**
		$C_{19}H_{13}F_3N_2O_2$	78863-62-4	
	InChI=1/C19H13F3N2O2/c20-19(21,22)13-6-4-9-15(12-13)26-18-16(10-5-11-23-18)17(25)24-14-7-2-1-3-8-14/h1-12H,(H,24,25)/f/h24H			

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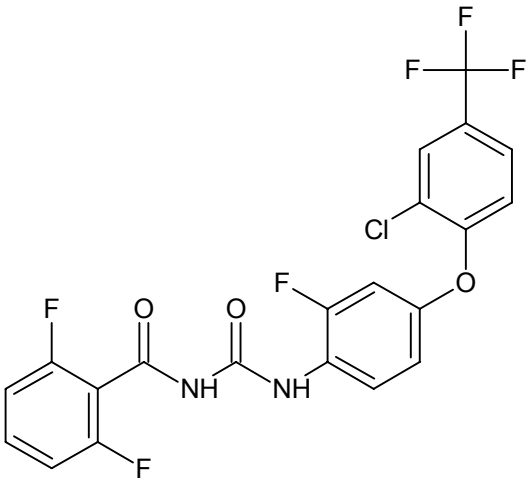
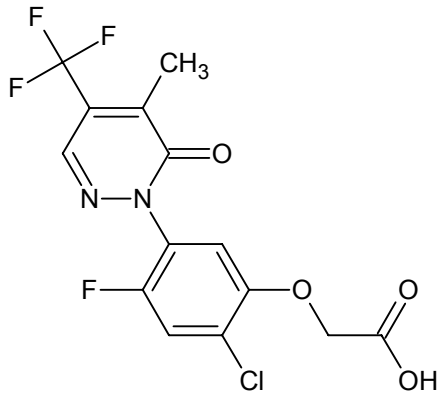
E flufenoxuron **F** flufénoxuron (m)	1-[4-(2-chloro-α,α,α-trifluoro-*p*-tolyloxy)-2-fluorophenyl]-3-(2,6-difluorobenzoyl)urea 1-[4-(2-chloro-4-trifluorométhylphénoxy)-2-fluorophényl]-3-(2,6-difluorobenzoyl)urée *N*-[[[4-[2-chloro-4-(trifluoromethyl)phenoxy]-2-fluorophenyl]amino]carbonyl]-2,6-difluorobenzamide			**A** **IGR**
		$C_{21}H_{11}ClF_6N_2O_3$	101463-69-8	
	InChI=1/C21H11ClF6N2O3/c22-12-8-10(21(26,27)28)4-7-17(12)33-11-5-6-16(15(25)9-11)29-20(32)30-19(31)18-13(23)2-1-3-14(18)24/h1-9H,(H2,29,30,31,32)/f/h29-30H			
E flufenpyr **F** flufenpyr (m)	2-chloro-5-[1,6-dihydro-5-methyl-6-oxo-4-(trifluoromethyl)pyridazin-1-yl]-4-fluorophenoxyacetic acid acide 2-chloro-5-[1,6-dihydro-5-méthyl-6-oxo-4-(trifluorométhyl)pyridazin-1-yl]-4-fluorophénoxyacétique [2-chloro-4-fluoro-5-[5-methyl-6-oxo-4-(trifluoromethyl)-1(6*H*)-pyridazinyl]phenoxy]acetic acid			**H**
		$C_{14}H_9ClF_4N_2O_4$	188490-07-5	
	InChI=1/C14H9ClF4N2O4/c1-6-7(14(17,18)19)4-20-21(13(6)24)10-3-11(25-5-12(22)23)8(15)2-9(10)16/h2-4H,5H2,1H3,(H,22,23)/f/h22H flufenpyr-ethyl InChI=1/C16H13ClF4N2O4/c1-3-26-14(24)7-27-13-5-12(11(18)4-10(13)17)23-15(25)8(2)9(6-22-23)16(19,20)21/h4-6H,3,7H2,1-2H3			
	NOTE It should be stated which ester or salt is present, for example flufenpyr-ethyl [188489-07-8]. NOTE Il convient de préciser quel est l'ester ou le sel présent, par exemple flufenpyr-éthyle [188489-07-8].			

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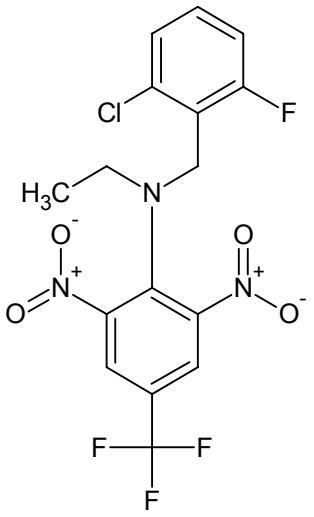
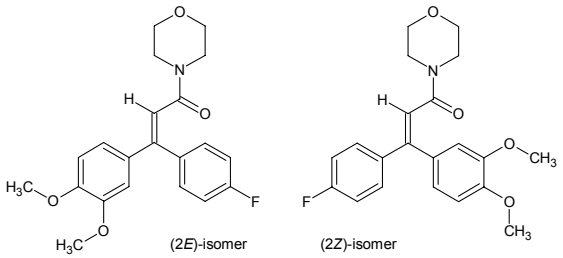
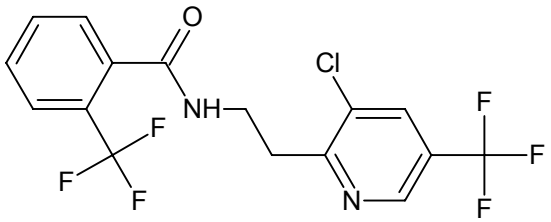
E flumetralin **F** flumétraline (f)	*N*-(2-chloro-6-fluorobenzyl)-*N*-ethyl-α,α,α-trifluoro-2,6-dinitro-*p*-toluidine *N*-(2-chloro-6-fluorobenzyl)-*N*-éthyl-4-trifluorométhyl-2,6-dinitroaniline 2-chloro-*N*-[2,6-dinitro-4-(trifluoromethyl)phenyl]-*N*-ethyl-6-fluorobenzenemethanamine			**P**
		$C_{16}H_{12}ClF_4N_3O_4$	62924-70-3	
	InChI=1/C16H12ClF4N3O4/c1-2-22(8-10-11(17)4-3-5-12(10)18)15-13(23(25)26)6-9(16(19,20)21)7-14(15)24(27)28/h3-7H,2,8H2,1H3			
E flumorph **F** flumorph (f)	(*EZ*)-3-(3,4-dimethoxyphenyl)-3-(4-fluorophenyl)-1-morpholinopropenone [50 % (*E*)-isomer, 50 % (*Z*)-isomer] (*EZ*)-3-(3,4-diméthoxyphényl)-3-(4-fluorophényl)-1-(morpholin-4-yl)acroléine [50 % d'isomère (*E*), 50 % d'isomère (*Z*)] 4-[3-(3,4-dimethoxyphenyl)-3-(4-fluorophenyl)-1-oxo-2-propenyl]morpholine			**F**
		$C_{21}H_{22}FNO_4$	211867-47-9	
	InChI=1/C21H22FNO4/c1-25-19-8-5-16(13-20(19)26-2)18(15-3-6-17(22)7-4-15)14-21(24)23-9-11-27-12-10-23/h3-8,13-14H,9-12H2,1-2H3/b18-14?			
E fluopyram **F** fluopyram (m)	*N*-{2-[3-chloro-5-(trifluoromethyl)-2-pyridyl]ethyl}-α,α,α-trifluoro-*o*-toluamide *N*-[2-[3-chloro-5-(trifluorométhyl))-2-pyridyl]éthyl]-α,α,α-trifluro-*o*-toluamide *N*-[2-[3-chloro-5-(trifluoromethyl)-2-pyridinyl]ethyl]-2-(trifluoromethyl)benzamide			**F**
		$C_{16}H_{11}ClF_6N_2O$	658066-35-4	
	InChI=1/C16H11ClF6N2O/c17-12-7-9(15(18,19)20)8-25-13(12)5-6-24-14(26)10-3-1-2-4-11(10)16(21,22)23/h1-4,7-8H,5-6H2,(H,24,26)/f/h24H			

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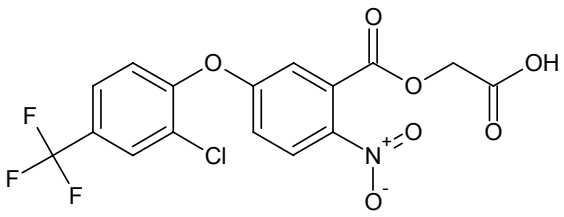
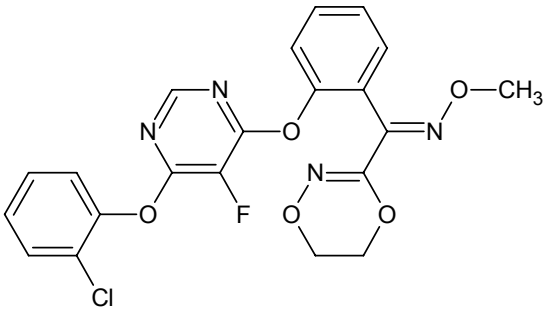
E fluoroglycofen **F** fluoroglycofène (m)	*O*-[5-(2-chloro-α,α,α-trifluoro-*p*-tolyloxy)-2-nitrobenzoyl]glycolic acid acide [5-(2-chloro-4-trifluorométhylphénoxy)-2-nitrobenzoyloxy]acétique carboxymethyl 5-[2-chloro-4-(trifluoromethyl)phenoxy]-2-nitrobenzoate	 $C_{16}H_9ClF_3NO_7$	77501-60-1	**H**
	InChI=1/C16H9ClF3NO7/c17-11-5-8(16(18,19)20)1-4-13(11)28-9-2-3-12(21(25)26)10(6-9)15(24)27-7-14(22)23/h1-6H,7H2,(H,22,23)/f/h22H fluoroglycofen-ethyl InChI=1/C18H13ClF3NO7/c1-2-28-16(24)9-29-17(25)12-8-11(4-5-14(12)23(26)27)30-15-6-3-10(7-13(15)19)18(20,21)22/h3-8H,2,9H2,1H3			
	NOTE It should be stated which ester or salt is present, for example fluoroglycofen-ethyl [77501-90-7]. NOTE Il convient de préciser quel est l'ester ou le sel présent, par exemple fluoroglycofène-éthyle [77501-90-7].			
E fluoxastrobin **F** fluoxastrobine (f)	(*E*)-{2-[6-(2-chlorophenoxy)-5-fluoropyrimidin-4-yloxy]phenyl}(5,6-dihydro-1,4,2-dioxazin-3-yl)methanone *O*-methyloxime (*E*)-{2-[6-(2-chlorophénoxy)-5-fluoropyrimidin-4-yloxy]phényl}-(5,6-dihydro-1,4,2-dioxazin-3-yl)méthanone-*O*-méthyloxime (1*E*)-[2-[[6-(2-chlorophenoxy)-5-fluoro-4-pyrimidinyl]oxy]phenyl](5,6-dihydro-1,4,2-dioxazin-3-yl)methanone *O*-methyloxime	 $C_{21}H_{16}ClFN_4O_5$	361377-29-9	**F**
	InChI=1/C21H16ClFN4O5/c1-28-26-18(21-27-30-11-10-29-21)13-6-2-4-8-15(13)31-19-17(23)20(25-12-24-19)32-16-9-5-3-7-14(16)22/h2-9,12H,10-11H2,1H3/b26-18+			

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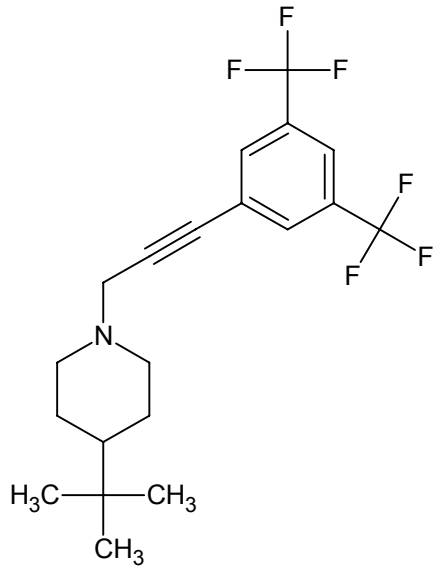
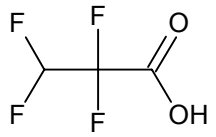
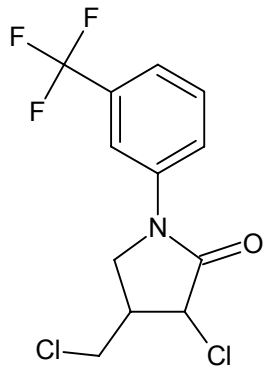
E flupropadine **F** flupropadine (f)	4-*tert*-butyl-1-[3-(α,α,α,α′,α′,α′-hexafluoro-3,5-xylyl)prop-2-ynyl]piperidine 4-*tert*-butyl-1-[1-(3,5-bis(trifluorométhyl)phényl)prop-2-ynyl]pipéridine 1-[3-[3,5-bis(trifluoromethyl)phenyl]-2-propynyl]-4-(1,1-dimethylethyl)piperidine			**R**
		$C_{20}H_{23}F_6N$	81613-59-4	
	InChI=1/C20H23F6N/c1-18(2,3)15-6-9-27(10-7-15)8-4-5-14-11-16(19(21,22)23)13-17(12-14)20(24,25)26/h11-13,15H,6-10H2,1-3H3			
E flupropanate **F** flupropanate (m)	2,2,3,3-tetrafluoropropionic acid acide 2,2,3,3-tétrafluoropropionique 2,2,3,3-tetrafluoropropanoic acid			**H**
		$C_3H_2F_4O_2$	756-09-2	
	InChI=1/C3H2F4O2/c4-1(5)3(6,7)2(8)9/h1H,(H,8,9)/f/h8H flupropanate-sodium InChI=1/C3H2F4O2.Na/c4-1(5)3(6,7)2(8)9;/h1H,(H,8,9);/q;+1/p-1/fC3HF4O2.Na/q-1;m			
	NOTE It should be stated which ester or salt is present, for example flupropanate-sodium [22898-01-7]. NOTE Il convient de préciser quel est l'ester ou le sel présent, par exemple flupropanate-sodium [22898-01-7].			
E flurochloridone **F** flurochloridone (f)	mixture of the enantiomeric pairs (3*RS*,4*RS*;3*RS*,4*SR*)-3-chloro-4-chloromethyl-1-(α,α,α-trifluoro-*m*-tolyl)-2-pyrrolidone (isomers in the ratio 3:1) mélange de paires d'énantiomères (3*RS*,4*RS*;3*RS*,4*SR*)-3-chloro-4-chlorométhyl-1-(3-trifluorométhylphényl)-pyrrolidone (isomères en proportion 3:1) 3-chloro-4-(chloromethyl)-1-[3-(trifluoromethyl)phenyl]-2-pyrrolidinone			**H**
		$C_{12}H_{10}Cl_2F_3NO$	61213-25-0	
	InChI=1/C12H10Cl2F3NO/c13-5-7-6-18(11(19)10(7)14)9-3-1-2-8(4-9)12(15,16)17/h1-4,7,10H,5-6H2/t7?,10?			

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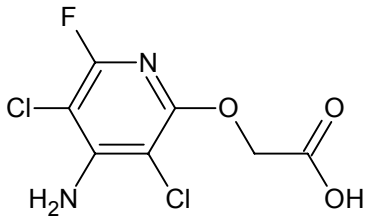
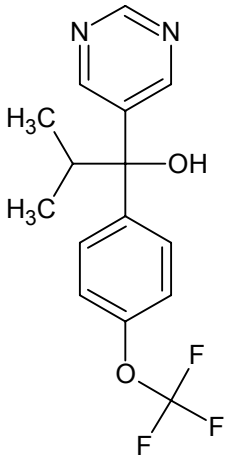
E fluroxypyr **F** fluoroxypyr (m)	4-amino-3,5-dichloro-6-fluoro-2-pyridyloxyacetic acid acide 4-amino-3,5-dichloro-6-fluoropyrid-2-yloxyacétique [(4-amino-3,5-dichloro-6-fluoro-2-pyridinyl)oxy]acetic acid			**H**
		$C_7H_5Cl_2FN_2O_3$	69377-81-7	
	InChI=1/C7H5Cl2FN2O3/c8-3-5(11)4(9)7(12-6(3)10)15-1-2(13)14/h1H2,(H2,11,12)(H,13,14)/f/h13H,11H2 fluroxypyr-butometyl InChI=1/C14H19Cl2FN2O4/c1-3-4-5-21-6-8(2)23-9(20)7-22-14-11(16)12(18)10(15)13(17)19-14/h8H,3-7H2,1-2H3,(H2,18,19)/t8-/s3/f/h18H2 fluroxypyr-meptyl InChI=1/C15H21Cl2FN2O3/c1-3-4-5-6-7-9(2)23-10(21)8-22-15-12(17)13(19)11(16)14(18)20-15/h9H,3-8H2,1-2H3,(H2,19,20)/t9-/s3/f/h19H2			
	NOTE It should be stated which ester or salt is present, for example fluroxypyr-butometyl [154486-27-8] or fluroxypyr-meptyl [81406-37-3]. NOTE Il convient de préciser quel est l'ester ou le sel présent, par exemple fluoroxypyr-butometyl [154486-27-8] ou fluoroxypyr-meptyl [81406-37-3].			
E flurprimidol **F** flurprimidol (m)	(*RS*)-2-methyl-1-pyrimidin-5-yl-1-(4-trifluoromethoxyphenyl)propan-1-ol (*RS*)-2-méthyl-1-(pyrimidin-5-yl)-1-(4-trifluorométhoxyphényl)propan-1-ol α-(1-methylethyl)-α-[4-(trifluoromethoxy)phenyl]-5-pyrimidinemethanol			**P**
		$C_{15}H_{15}F_3N_2O_2$	56425-91-3	
	InChI=1/C15H15F3N2O2/c1-10(2)14(21,12-7-19-9-20-8-12)11-3-5-13(6-4-11)22-15(16,17)18/h3-10,21H,1-2H3/t14-/s3			

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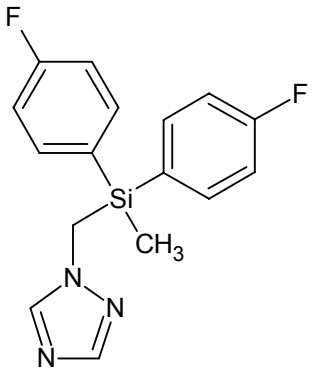
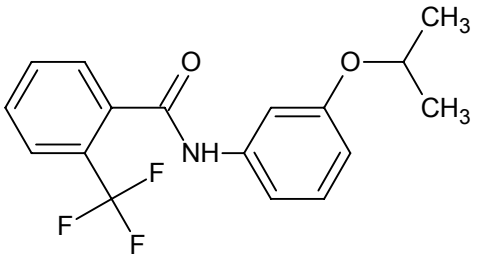
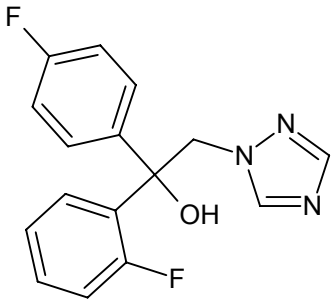
E flusilazole **F** flusilazole (m)	bis(4-fluorophenyl)(methyl)(1*H*-1,2,4-triazol-1-ylmethyl)silane 1-[[bis(4-fluorophényl)méthylsilyl]méthyl]-1*H*-1,2,4-triazole 1-[[bis(4-fluorophenyl)methylsilyl]methyl]-1*H*-1,2,4-triazole			**F**
		$C_{16}H_{15}F_2N_3Si$	85509-19-9	
	InChI=1/C16H15F2N3Si/c1-22(12-21-11-19-10-20-21,15-6-2-13(17)3-7-15)16-8-4-14(18)5-9-16/h2-11H,12H2,1H3			
E flutolanil **F** flutolanil (m)	α,α,α-trifluoro-3′-isopropoxy-*o*-toluanilide 3-isopropoxy-2′-trifluorométhylbenzanilide *N*-[3-(1-methylethoxy)phenyl]-2-(trifluoromethyl)benzamide			**F**
		$C_{17}H_{16}F_3NO_2$	66332-96-5	
	InChI=1/C17H16F3NO2/c1-11(2)23-13-7-5-6-12(10-13)21-16(22)14-8-3-4-9-15(14)17(18,19)20/h3-11H,1-2H3,(H,21,22)/f/h21H			
	NOTE The name “flutolanil” is not acceptable for use in Hungary because of the risk of confusion with the trade name “Flutolan”. NOTE Le nom «flutolanil» n'est pas acceptable pour l'emploi en Hongrie, car il entre en conflit avec le nom commercial «Flutolan».			
E flutriafol **F** flutriafol (m)	(*RS*)-2,4′-difluoro-α-(1*H*-1,2,4-triazol-1-ylmethyl)benzhydryl alcohol (*RS*)-1-(2-fluorophényl)-1-(4-fluorophényl)-2-(1*H*-1,2,4-triazol-1-ylméthyl)éthanol α-(2-fluorophenyl)-α-(4-fluorophenyl)-1*H*-1,2,4-triazole-1-ethanol			**F**
		$C_{16}H_{13}F_2N_3O$	76674-21-0	
	InChI=1/C16H13F2N3O/c17-13-7-5-12(6-8-13)16(22,9-21-11-19-10-20-21)14-3-1-2-4-15(14)18/h1-8,10-11,22H,9H2/t16-/s3			

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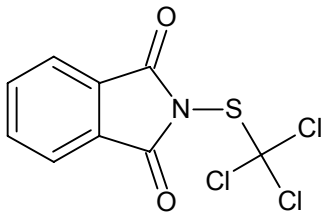
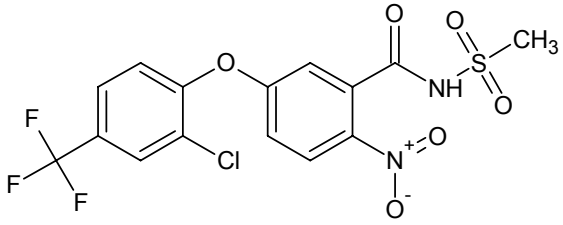
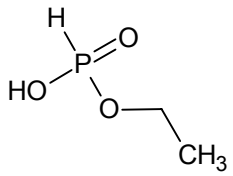
E folpet **F** folpet (m)	*N*-(trichloromethylthio)phthalimide *N*-(trichlorométhylthio)phtalimide 2-[(trichloromethyl)thio]-1*H*-isoindole-1,3(2*H*)-dione			**F**
		$C_9H_4Cl_3NO_2S$	133-07-3	
	InChI=1/C9H4Cl3NO2S/c10-9(11,12)16-13-7(14)5-3-1-2-4-6(5)8(13)15/h1-4H			
	NOTE The name “folpet” is not acceptable for use in France, where “folpel” has been adopted as the common name. NOTE Le nom «folpet» n'est pas acceptable pour l'emploi en France, où «folpel» a été adopté comme nom commun.			
E fomesafen **F** fomésafène (m)	5-(2-chloro-α,α,α-trifluoro-*p*-tolyloxy)-*N*-mesyl-2-nitrobenzamide 5-(2-chloro-4-trifluorométhylphénoxy)-*N*-mésyl-2-nitrobenzamide 5-[2-chloro-4-(trifluoromethyl)phenoxy]-*N*-(methylsulfonyl)-2-nitrobenzamide			**H**
		$C_{15}H_{10}ClF_3N_2O_6S$	72178-02-0	
	InChI=1/C15H10ClF3N2O6S/c1-28(25,26)20-14(22)10-7-9(3-4-12(10)21(23)24)27-13-5-2-8(6-11(13)16)15(17,18)19/h2-7H,1H3,(H,20,22)/f/h20H fomesafen-sodium InChI=1/C15H10ClF3N2O6S.Na/c1-28(25,26)20-14(22)10-7-9(3-4-12(10)21(23)24)27-13-5-2-8(6-11(13)16)15(17,18)19;/h2-7H,1H3,(H,20,22);/q;+1/p-1/fC15H9ClF3N2O6S.Na/q-1;m/b20-14?;			
	NOTE It should be stated which salt is present, for example fomesafen-sodium [108731-70-0]. NOTE Il convient de préciser quel est le sel présent, par exemple fomésafène-sodium [108731-70-0].			
E fosetyl **F** fosétyl (m)	ethyl hydrogen phosphonate hydrogénophosphonate d'éthyle ethyl hydrogen phosphonate			**F**
		$C_2H_7O_3P$	15845-66-6	
	InChI=1/C2H7O3P/c1-2-5-6(3)4/h6H,2H2,1H3,(H,3,4)/f/h3H fosetyl-aluminium InChI=1/3C2H7O3P.Al/c3*1-2-5-6(3)4;/h3*6H,2H2,1H3,(H,3,4);/q;;;+3/p-3/f3C2H6O3P.Al/q3*-1;m			
	NOTE It should be stated which ester or salt is present, for example fosetyl-aluminium [39148-24-8]. NOTE Il convient de préciser quel est l'ester ou le sel présent, par exemple fosétyl-aluminium [39148-24-8].			

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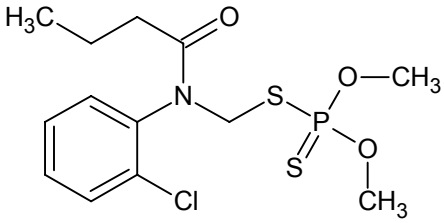
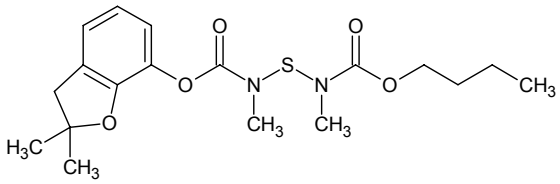
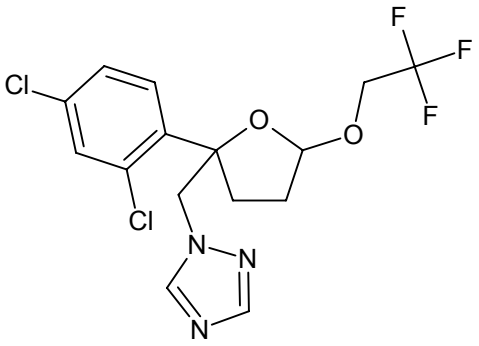
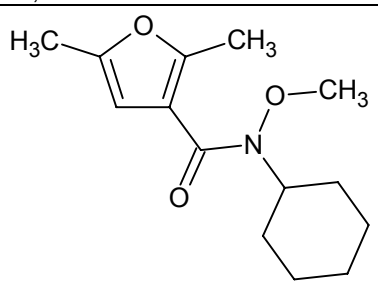
E fosmethilan **F** fosméthilane (m)	2′-chloro-*N*-(dimethoxy= phosphinothioylthiomethyl)butyranilide dithiophosphate de *S*-(butyryl-2-chlorophényl-amino)méthyle et de *O,O*-diméthyle *S*-[[(2-chlorophenyl)(1-oxobutyl)amino]methyl] *O,O*-dimethyl phosphorodithioate			**I**
		$C_{13}H_{19}ClNO_3PS_2$	83733-82-8	
	InChI=1/C13H19ClNO3PS2/c1-4-7-13(16)15(10-21-19(20,17-2)18-3)12-9-6-5-8-11(12)14/h5-6,8-9H,4,7,10H2,1-3H3			
E furathiocarb **F** furathiocarb (m)	butyl 2,3-dihydro-2,2-dimethylbenzofuran-7-yl *N,N′*-dimethyl-*N,N′*-thiodicarbamate thiobis(méthylcarbamate) de butyle et de 2,2-diméthyl-2,3-dihydrobenzofuran-7-yle 2,3-dihydro-2,2-dimethyl-7-benzofuranyl 2,4-dimethyl-5-oxo-6-oxa-3-thia-2,4-diazadecanoate			**I**
		$C_{18}H_{26}N_2O_5S$	65907-30-4	
	InChI=1/C18H26N2O5S/c1-6-7-11-23-16(21)19(4)26-20(5)17(22)24-14-10-8-9-13-12-18(2,3)25-15(13)14/h8-10H,6-7,11-12H2,1-5H3			
E furconazole **F** furconazole (m)	(2*RS*,5*RS*;2*RS*,5*SR*)-5-(2,4-dichlorophenyl)tetrahydro-5-(1*H*-1,2,4-triazol-1-ylmethyl)-2-furyl 2,2,2-trifluoroethyl ether (2*RS*,5*RS*;2*RS*,5*SR*)-1-[[2-(2,4-dichlorophényl)-5-(2,2,2-trifluoroéthoxy)-tétrahydrofuran-2-yl]méthyl]-1*H*-1,2,4-triazole 1-[[2-(2,4-dichlorophenyl)tetrahydro-5-(2,2,2-trifluoroethoxy)-2-furanyl]methyl]-1*H*-1,2,4-triazole			**F**
		$C_{15}H_{14}Cl_2F_3N_3O_2$	112839-33-5	
	InChI=1/C15H14Cl2F3N3O2/c16-10-1-2-11(12(17)5-10)14(6-23-9-21-8-22-23)4-3-13(25-14)24-7-15(18,19)20/h1-2,5,8-9,13H,3-4,6-7H2/t13?,14?			
E furmecyclox **F** furmécyclox (m)	methyl *N*-cyclohexyl-2,5-dimethyl-3-furohydroxamate *N*-cyclohexyl-*N*-méthoxy-2,5-diméthylfuramide *N*-cyclohexyl-*N*-methoxy-2,5-dimethyl-3-furancarboxamide			**F**
		$C_{14}H_{21}NO_3$	60568-05-0	
	InChI=1/C14H21NO3/c1-10-9-13(11(2)18-10)14(16)15(17-3)12-7-5-4-6-8-12/h9,12H,4-8H2,1-3H3			

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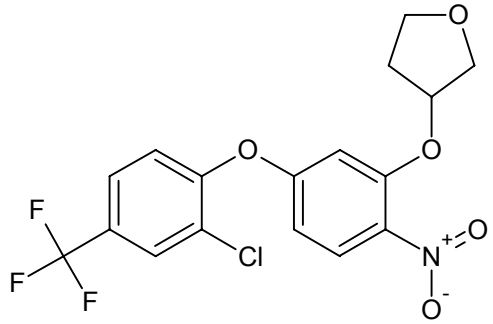
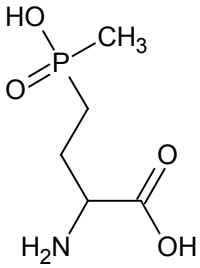
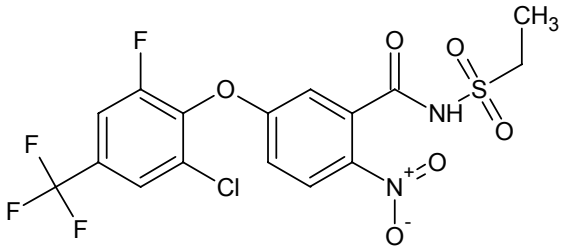
E furyloxyfen **F** furyloxyfène (m)	(*RS*)-5-(2-chloro-α,α,α-trifluoro-*p*-tolyloxy)-2-nitrophenyl tetrahydro-3-furyl ether oxyde de 5-(2-chloro-4-trifluorométhyl)phénoxy-2-nitrophényle et de (*RS*)-fur-3-yle 3-[5-[2-chloro-4-(trifluoromethyl)phenoxy]-2-nitrophenoxy]tetrahydrofuran			**H**
		$C_{17}H_{13}ClF_3NO_5$	80020-41-3	
	InChI=1/C17H13ClF3NO5/c18-13-7-10(17(19,20)21)1-4-15(13)26-11-2-3-14(22(23)24)16(8-11)27-12-5-6-25-9-12/h1-4,7-8,12H,5-6,9H2/t12-/s3			
E glufosinate **F** glufosinate (m)	(*RS*)-2-amino-4-[hydroxy(methyl)=phosphinoyl]butyric acid acide (*RS*)-4-[hydroxy(méthyl)phosphinoyl]-2-aminobutyrique 2-amino-4-(hydroxymethylphosphinyl)butanoic acid			**H**
		$C_5H_{12}NO_4P$	51276-47-2	
	InChI=1/C5H12NO4P/c1-11(9,10)3-2-4(6)5(7)8/h4H,2-3,6H2,1H3,(H,7,8)(H,9,10)/t4-/s3/f/h7,9H glufosinate-ammonium InChI=1/C5H12NO4P.H3N/c1-11(9,10)3-2-4(6)5(7)8;/h4H,2-3,6H2,1H3,(H,7,8)(H,9,10);1H3/t4-;/s3/fC5H11NO4P.H4N/h9H;1H/q-1;+1			
	NOTE It should be stated which ester or salt is present, for example glufosinate-ammonium [77182-82-2]. NOTE Il convient de préciser quel est l'ester ou le sel présent, par exemple glufosinate-ammonium [77182-82-2].			
E halosafen **F** halosafène (m)	5-(2-chloro-α,α,α,6-tetrafluoro-*p*-tolyloxy)-*N*-ethylsulfonyl-2-nitrobenzamide 5-(2-chloro-6-fluoro-4-trifluorométhylphénoxy)-*N*-éthylsulfonyl-2-nitrobenzamide 5-[2-chloro-6-fluoro-4-(trifluoromethyl)phenoxy]-*N*-(ethylsulfonyl)-2-nitrobenzamide			**H**
		$C_{16}H_{11}ClF_4N_2O_6S$	77227-69-1	
	InChI=1/C16H11ClF4N2O6S/c1-2-30(27,28)22-15(24)10-7-9(3-4-13(10)23(25)26)29-14-11(17)5-8(6-12(14)18)16(19,20)21/h3-7H,2H2,1H3,(H,22,24)/f/h22H			

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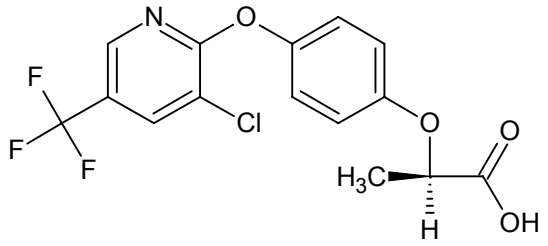
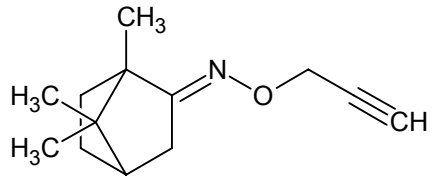
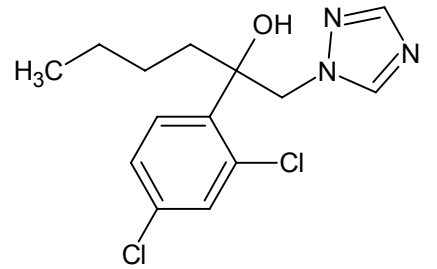
E haloxyfop-P **F** haloxyfop-P (m)	(*R*)-2-{4-[3-chloro-5-(trifluoromethyl)-2-pyridyloxy]phenoxy}propionic acid acide (*R*)-2-{4-[3-chloro-5-(trifluorométhyl)pyrid-2-yloxy]phénoxy}propionique (2*R*)-2-[4-[[3-chloro-5-(trifluoromethyl)-2-pyridinyl]oxy]phenoxy]propanoic acid			**H**
		$C_{15}H_{11}ClF_3NO_4$	95977-29-0	
	InChI=1/C15H11ClF3NO4/c1-8(14(21)22)23-10-2-4-11(5-3-10)24-13-12(16)6-9(7-20-13)15(17,18)19/h2-8H,1H3,(H,21,22)/t8-/m1/s1/f/h21H haloxyfop-P-methyl InChI=1/C16H13ClF3NO4/c1-9(15(22)23-2)24-11-3-5-12(6-4-11)25-14-13(17)7-10(8-21-14)16(18,19)20/h3-9H,1-2H3/t9-/m1/s1			
	NOTE It should be stated which ester or salt is present, for example haloxyfop-P-methyl [72619-32-0]. NOTE Il convient de préciser quel est l'ester ou le sel présent, par exemple haloxyfop-P-méthyl [72619-32-0].			
E heptopargil **F** heptopargil (m)	(1*RS*,4*RS*)-bornan-2-one (*E*)-*O*-prop-2-ynyloxime (1*RS*,4*RS*)-bornan-2-one-(*E*)-*O*-prop-2-ynyloxime (1*R*,2*E*,4*R*)-*rel*-1,7,7-trimethylbicyclo[2.2.1]heptan-2-one *O*-2-propynyloxime			**P**
		$C_{13}H_{19}NO$	73886-28-9	
	InChI=1/C13H19NO/c1-5-8-15-14-11-9-10-6-7-13(11,4)12(10,2)3/h1,10H,6-9H2,2-4H3/b14-11+/t10-,13+/s3			
E hexaconazole **F** hexaconazole (m)	(*RS*)-2-(2,4-dichlorophenyl)-1-(1*H*-1,2,4-triazol-1-yl)hexan-2-ol (*RS*)-2-(2,4-dichlorophényl)-1-(1*H*-1,2,4-triazol-1-yl)hexan-2-ol α-butyl-α-(2,4-dichlorophenyl)-1*H*-1,2,4-triazole-1-ethanol			**F**
		$C_{14}H_{17}Cl_2N_3O$	79983-71-4	
	InChI=1/C14H17Cl2N3O/c1-2-3-6-14(20,8-19-10-17-9-18-19)12-5-4-11(15)7-13(12)16/h4-5,7,9-10,20H,2-3,6,8H2,1H3/t14-/s3			

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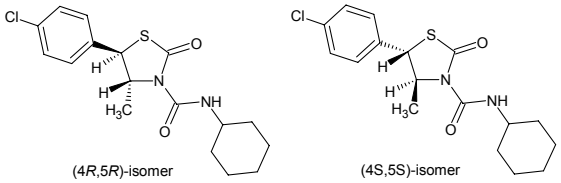
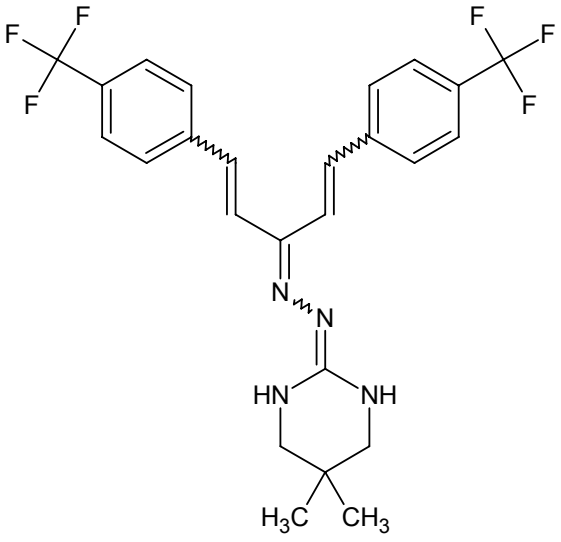
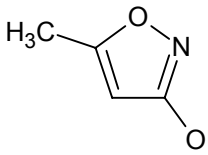
E hexythiazox **F** hexythiazox (m)	(4*RS*,5*RS*)-5-(4-chlorophenyl)-*N*-cyclohexyl-4-methyl-2-oxo-1,3-thiazolidine-3-carboxamide (4*RS*,5*RS*)-5-(4-chlorophényl)-*N*-cyclohexyl-4-méthyl-2-oxo-1,3-thiazolidine-3-carboxamide (4*R*,5*R*)-*rel*-5-(4-chlorophenyl)-*N*-cyclohexyl-4-methyl-2-oxo-3-thiazolidinecarboxamide	 (4*R*,5*R*)-isomer (4S,5S)-isomer $C_{17}H_{21}ClN_2O_2S$	78587-05-0	**A**
	InChI=1/C17H21ClN2O2S/c1-11-15(12-7-9-13(18)10-8-12)23-17(22)20(11)16(21)19-14-5-3-2-4-6-14/h7-11,14-15H,2-6H2,1H3,(H,19,21)/t11-,15+/s3/f/h19H			
E hydramethylnon **F** hydraméthylnone (f)	5,5-dimethylperhydropyrimidin-2-one 4-trifluoromethyl-α-(4-trifluoromethylstyryl)= cinnamylidenehydrazone 5,5-diméthyl-tétrahydro-1*H*-pyrimidin-2-one-bis-(4-trifluorométhylstyryl)-hydrazone tetrahydro-5,5-dimethyl-2(1*H*)-pyrimidinone [3-[4-(trifluoromethyl)phenyl]-1-[2-[4-(trifluoromethyl)phenyl]ethenyl]-2-propenylidene]hydrazone	 $C_{25}H_{24}F_6N_4$	67485-29-4	**I**
	InChI=1/C25H24F6N4/c1-23(2)15-32-22(33-16-23)35-34-21(13-7-17-3-9-19(10-4-17)24(26,27)28)14-8-18-5-11-20(12-6-18)25(29,30)31/h3-14H,15-16H2,1-2H3,(H2,32,33,35)/b13-7?,14-8?,34-21?/f/h32-33H			
E hymexazol **F** hymexazol (m)	5-methyl-1,2-oxazol-3-ol 5-méthyl-isoxazol-3-ol 5-methyl-3(2*H*)-isoxazolone	 $C_4H_5NO_2$	10004-44-1	**F** **P**
	InChI=1/C4H5NO2/c1-3-2-4(6)5-7-3/h2H,1H3,(H,5,6)/f/h6H CAS-preferred tautomer InChI=1/C4H5NO2/c1-3-2-4(6)5-7-3/h2H,1H3,(H,5,6)/f/h5H			
	NOTE The CAS name and Registry Number are for the tautomer that is preferred under the CAS nomenclature rules. NOTE Le nom et le numéro du CAS sont pour le tautomère, qui est la forme préférée selon les règles de nomenclature CAS.			

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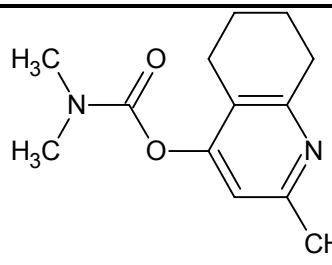
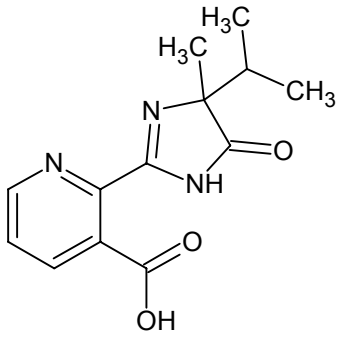
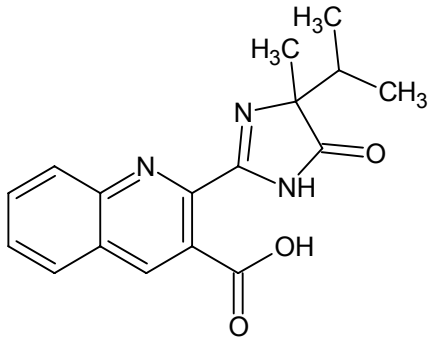
E hyquincarb **F** hyquincarb (m)	5,6,7,8-tetrahydro-2-methyl-4-quinolyl dimethylcarbamate diméthylcarbamate de 5,6,7,8-tétrahydro-2-méthylquinol-4-yle 5,6,7,8-tetrahydro-2-methyl-4-quinolinyl dimethylcarbamate			**I**
		$C_{13}H_{18}N_2O_2$	56716-21-3	
	InChI=1/C13H18N2O2/c1-9-8-12(17-13(16)15(2)3)10-6-4-5-7-11(10)14-9/h8H,4-7H2,1-3H3			
E imazapyr **F** imazapyr (m)	(*RS*)-2-(4-isopropyl-4-methyl-5-oxo-2-imidazolin-2-yl)nicotinic acid acide (*RS*)-2-(4-isopropyl-4-méthyl-5-oxo-imidazolin-2-yl)nicotinique 2-[4,5-dihydro-4-methyl-4-(1-methylethyl)-5-oxo-1*H*-imidazol-2-yl]-3-pyridinecarboxylic acid			**H**
		$C_{13}H_{15}N_3O_3$	81334-34-1	
	InChI=1/C13H15N3O3/c1-7(2)13(3)12(19)15-10(16-13)9-8(11(17)18)5-4-6-14-9/h4-7H,1-3H3,(H,17,18)(H,15,16,19)/t13-/s3/f/h15,17H imazapyr-isopropylammonium InChI=1/C13H15N3O3.C3H9N/c1-7(2)13(3)12(19)15-10(16-13)9-8(11(17)18)5-4-6-14-9;1-3(2)4/h4-7H,1-3H3,(H,17,18)(H,15,16,19);3H,4H2,1-2H3/t13-;/s3/f/h15,17H;			
	NOTE It should be stated which ester or salt is present, for example imazapyr-isopropylammonium [81510-83-0]. NOTE Il convient de préciser quel est l'ester ou le sel présent, par exemple imazapyr-isopropylammonium [81510-83-0].			
E imazaquin **F** imazaquine (m)	(*RS*)-2-(4-isopropyl-4-methyl-5-oxo-2-imidazolin-2-yl)quinoline-3-carboxylic acid acide (*RS*)-2-(4-isopropyl-4-méthyl-5-oxo-2-imidazolin-2-yl)quinoline-3-carboxylique 2-[4,5-dihydro-4-methyl-4-(1-methylethyl)-5-oxo-1*H*-imidazol-2-yl]-3-quinolinecarboxylic acid			**H**
		$C_{17}H_{17}N_3O_3$	81335-37-7	
	InChI=1/C17H17N3O3/c1-9(2)17(3)16(23)19-14(20-17)13-11(15(21)22)8-10-6-4-5-7-12(10)18-13/h4-9H,1-3H3,(H,21,22)(H,19,20,23)/t17-/s3/f/h19,21H imazaquin-ammonium InChI=1/C17H17N3O3.H3N/c1-9(2)17(3)16(23)19-14(20-17)13-11(15(21)22)8-10-6-4-5-7-12(10)18-13;/h4-9H,1-3H3,(H,21,22)(H,19,20,23);1H3/t17-;/s3/fC17H16N3O3.H4N/h19H;1H/q-1;+1			
	NOTE It should be stated which ester or salt is present, for example imazaquin-ammonium [81335-47-9]. NOTE Il convient de préciser quel est l'ester ou le sel présent, par exemple imazaquine-ammonium [81335-47-9].			

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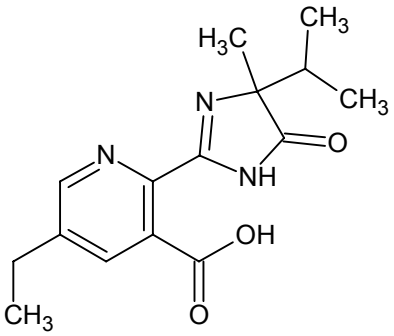
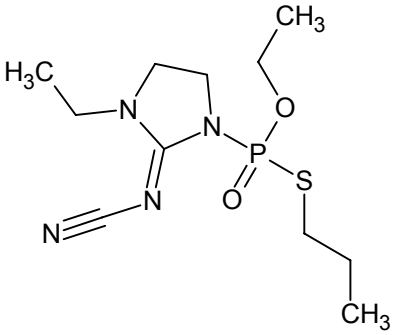
E imazethapyr **F** imazéthapyr (m)	(*RS*)-5-ethyl-2-(4-isopropyl-4-methyl-5-oxo-2-imidazolin-2-yl)nicotinic acid acide (*RS*)-5-éthyl-2-(4-isopropyl-4-méthyl-5-oxo-imidazolin-2-yl)nicotinique 2-[4,5-dihydro-4-methyl-4-(1-methylethyl)-5-oxo-1*H*-imidazol-2-yl]-5-ethyl-3-pyridinecarboxylic acid			**H**
		$C_{15}H_{19}N_3O_3$	81335-77-5	
	InChI=1/C15H19N3O3/c1-5-9-6-10(13(19)20)11(16-7-9)12-17-14(21)15(4,18-12)8(2)3/h6-8H,5H2,1-4H3,(H,19,20)(H,17,18,21)/t15-/s3/f/h17,19H imazethapyr-ammonium InChI=1/C15H19N3O3.H3N/c1-5-9-6-10(13(19)20)11(16-7-9)12-17-14(21)15(4,18-12)8(2)3;/h6-8H,5H2,1-4H3,(H,19,20)(H,17,18,21);1H3/t15-;/s3/fC15H18N3O3.H4N/h17H;1H/q-1;+1			
	NOTE It should be stated which ester or salt is present, for example imazethapyr-ammonium [101917-66-2]. NOTE Il convient de préciser quel est l'ester ou le sel présent, par exemple imazéthapyr-ammonium [101917-66-2].			
E imicyafos **F** imicyafos (m)	(*RS*)-{*O*-ethyl *S*-propyl (*E*)-[2-(cyanoimino)-3-ethylimidazolidin-1-yl]phosphonothioate} (*RS*)-{(*E*)-[2-(cyanoimino)-3-éthylimidazolidin-1-yl]phosphonothioate de *O*-éthyle et de *S*-propyle} *O*-ethyl *S*-propyl [(2*E*)-2-(cyanoimino)-3-ethyl-1-imidazolidinyl]phosphonothioate			**N**
		$C_{11}H_{21}N_4O_2PS$	140163-89-9	
	InChI=1/C11H21N4O2PS/c1-4-9-19-18(16,17-6-3)15-8-7-14(5-2)11(15)13-10-12/h4-9H2,1-3H3/b13-11+/t18-/s3			

E: Common name F: Nom commun	Chemical name Nom chimique E: IUPAC F: UICPA C: CAS	Structure Structure Molecular formula Formule brute	CAS Registry Number® Numéro d'enregistrement CAS®	Use Application
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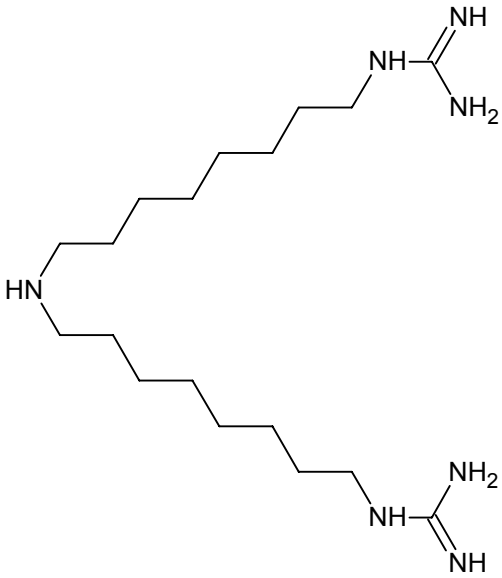
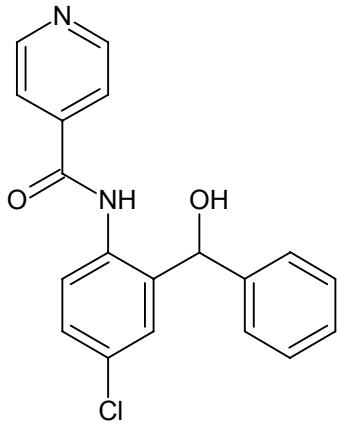
E iminoctadine **F** iminoctadine (f)	1,1′-(iminodioctamethylene)diguanidine bis(8-guanidinoctyl)amine ou 1,1'-(iminodioctaméthylène)diguanidine *N,N‴*-(iminodi-8,1-octanediyl)bis[guanidine]			**F**
		$C_{18}H_{41}N_7$	13516-27-3	
	InChI=1/C18H41N7/c19-17(20)24-15-11-7-3-1-5-9-13-23-14-10-6-2-4-8-12-16-25-18(21)22/h23H,1-16H2,(H4,19,20,24)(H4,21,22,25)/f/h19,21,24-25H,20,22H2 iminoctadine triacetate InChI=1/C18H41N7.3C2H4O2/c19-17(20)24-15-11-7-3-1-5-9-13-23-14-10-6-2-4-8-12-16-25-18(21)22;3*1-2(3)4/h23H,1-16H2,(H4,19,20,24)(H4,21,22,25);3*1H3,(H,3,4)/f/h19,21,24-25H,20,22H2;3*3H			
	NOTE 1 It should be stated which salt is present, for example iminoctadine triacetate [57520-17-9]. NOTE 1 Il convient de préciser quel est le sel présent, par exemple iminoctadine triacétate [57520-17-9]. NOTE 2 The name is a replacement for the common name “guazatine”, the definition of which has been changed. NOTE 2 Le nom remplace le nom commun «guazatine», dont la définition a été modifiée.			
E inabenfide **F** inabenfide (m)	(*RS*)-4′-chloro-2′-(α-hydroxybenzyl)isonicotinanilide (*RS*)-4′-chloro-2′-(hydroxy(phényl)=méthyl)isonicotinanilide *N*-[4-chloro-2-(hydroxyphenylmethyl)phenyl]-4-pyridinecarboxamide			**P**
		$C_{19}H_{15}ClN_2O_2$	82211-24-3	
	InChI=1/C19H15ClN2O2/c20-15-6-7-17(22-19(24)14-8-10-21-11-9-14)16(12-15)18(23)13-4-2-1-3-5-13/h1-12,18,23H,(H,22,24)/t18-/s3/f/h22H			

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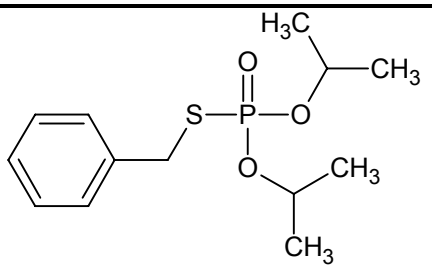
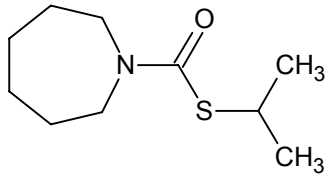
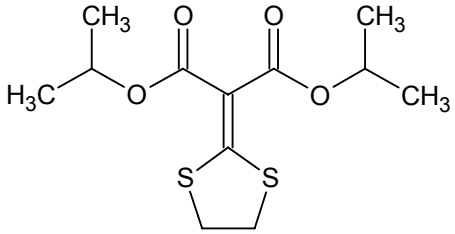
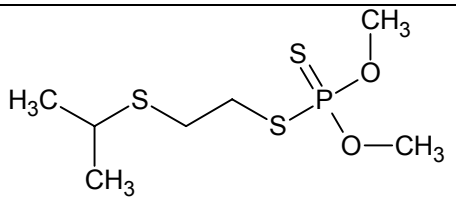
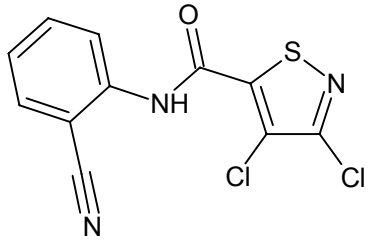
E iprobenfos **F** iprobenfos (m)	*S*-benzyl *O,O*-diisopropyl phosphorothioate thiophosphate de *S*-benzyle et de *O,O*-diisopropyle *O,O*-bis(1-methylethyl) *S*-(phenylmethyl) phosphorothioate			**F** **Y**
		$C_{13}H_{21}O_3PS$	26087-47-8	
	InChI=1/C13H21O3PS/c1-11(2)15-17(14,16-12(3)4)18-10-13-8-6-5-7-9-13/h5-9,11-12H,10H2,1-4H3			
E isopolinate **F** isopolinate (m)	*S*-isopropyl azepane-1-carbothioate hexahydro-1*H*-azépine-1-thiocarboxylate de *S*-isopropyle *S*-(1-methylethyl) hexahydro-1*H*-azepine-1-carbothioate			**H**
		$C_{10}H_{19}NOS$	3134-70-1	
	InChI=1/C10H19NOS/c1-9(2)13-10(12)11-7-5-3-4-6-8-11/h9H,3-8H2,1-2H3			
E isoprothiolane **F** isoprothiolane (m)	diisopropyl 1,3-dithiolan-2-ylidenemalonate 1,3-dithiolan-2-ylidènemalonate de diisopropyle bis(1-methylethyl) 1,3-dithiolan-2-ylidenepropanedioate			**F** **I**
		$C_{12}H_{18}O_4S_2$	50512-35-1	
	InChI=1/C12H18O4S2/c1-7(2)15-10(13)9(11(14)16-8(3)4)12-17-5-6-18-12/h7-8H,5-6H2,1-4H3			
E isothioate **F** isothioate (m)	*S*-2-isopropylthioethyl *O,O*-dimethyl phosphorodithioate phosphorodithioate de *S*-(2-isopropylthioéthyle) et de *O,O*-diméthyle *O,O*-dimethyl *S*-[2-[(1-methylethyl)thio]ethyl] phosphorodithioate			**I**
		$C_7H_{17}O_2PS_3$	36614-38-7	
	InChI=1/C7H17O2PS3/c1-7(2)12-5-6-13-10(11,8-3)9-4/h7H,5-6H2,1-4H3			
E isotianil **F** isotianil (m)	3,4-dichloro-2′-cyano-1,2-thiazole-5-carboxanilide *N*-(2-cyanophényl)-3,4-dichloroisothiazol-5-carboxamide 3,4-dichloro-*N*-(2-cyanophenyl)-5-isothiazolecarboxamide			**F**
		$C_{11}H_5Cl_2N_3OS$	224049-04-1	
	InChI=1/C11H5Cl2N3OS/c12-8-9(18-16-10(8)13)11(17)15-7-4-2-1-3-6(7)5-14/h1-4H,(H,15,17)/f/h15H			

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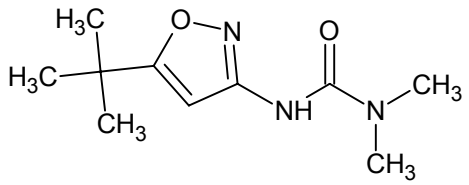
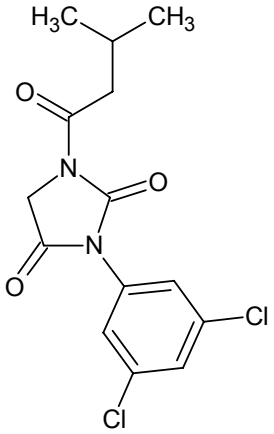
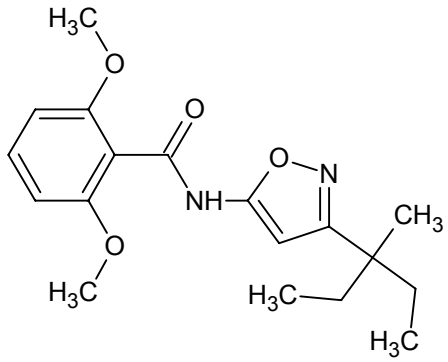
E isouron **F** isuron (m)	3-(5-*tert*-butyl-1,2-oxazol-3-yl)-1,1-dimethylurea *N*-(5-*tert*-butylisoxazol-3-yl)-*N′,N′*-diméthylurée ou 3-(5-*tert*-butylisoxazol-3-yl)-1,1-diméthylurée *N′*-[5-(1,1-dimethylethyl)-3-isoxazolyl]-*N,N*-dimethylurea			**H**
		$C_{10}H_{17}N_3O_2$	55861-78-4	
	InChI=1/C10H17N3O2/c1-10(2,3)7-6-8(12-15-7)11-9(14)13(4)5/h6H,1-5H3,(H,11,12,14)/f/h11H			
E isovaledione **F** isovalédione (f)	3-(3,5-dichlorophenyl)-1-isovalerylhydantoin 3-(3,5-dichlorophényl)-1-isovalérylhydantoïne 3-(3,5-dichlorophenyl)-1-(3-methyl-1-oxobutyl)-2,4-imidazolidinedione			**F**
		$C_{14}H_{14}Cl_2N_2O_3$	70017-93-5	
	InChI=1/C14H14Cl2N2O3/c1-8(2)3-12(19)17-7-13(20)18(14(17)21)11-5-9(15)4-10(16)6-11/h4-6,8H,3,7H2,1-2H3			
E isoxaben **F** isoxabène (m)	*N*-[3-(1-ethyl-1-methylpropyl)-1,2-oxazol-5-yl]-2,6-dimethoxybenzamide *N*-[3-(3-méthylpent-3-yl)isoxazol-5-yl]-2,6-diméthoxybenzamide *N*-[3-(1-ethyl-1-methylpropyl)-5-isoxazolyl]-2,6-dimethoxybenzamide			**H**
		$C_{18}H_{24}N_2O_4$	82558-50-7	
	InChI=1/C18H24N2O4/c1-6-18(3,7-2)14-11-15(24-20-14)19-17(21)16-12(22-4)9-8-10-13(16)23-5/h8-11H,6-7H2,1-5H3,(H,19,21)/f/h19H			

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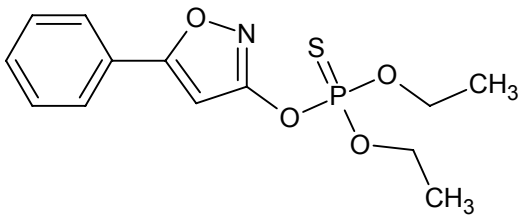
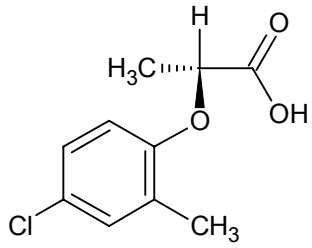
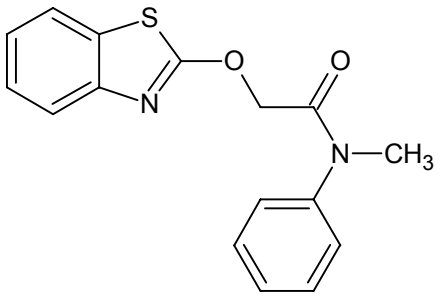
	IUPAC International Chemical Identifier (InChI™)			
E isoxathion **F** isoxathion (m)	*O,O*-diethyl *O*-5-phenyl-1,2-oxazol-3-yl phosphorothioate phosphorothioate de *O,O*-diéthyle et de *O*-(5-phényl-isoxazol-3-yle) *O,O*-diethyl *O*-(5-phenyl-3-isoxazolyl) phosphorothioate	 $C_{13}H_{16}NO_4PS$	18854-01-8	**I**
	InChI=1/C13H16NO4PS/c1-3-15-19(20,16-4-2)18-13-10-12(17-14-13)11-8-6-5-7-9-11/h5-10H,3-4H2,1-2H3			
E mecoprop-P **F** mécoprop-P (m)	(*R*)-2-(4-chloro-*o*-tolyloxy)propionic acid acide (*R*)-2-(4-chloro-2-méthyl-phénoxy)-propionique (2*R*)-2-(4-chloro-2-methylphenoxy)propanoic acid	 $C_{10}H_{11}ClO_3$	16484-77-8	**H**
	InChI=1/C10H11ClO3/c1-6-5-8(11)3-4-9(6)14-7(2)10(12)13/h3-5,7H,1-2H3,(H,12,13)/t7-/m1/s1/f/h12H mecoprop-P-isobutyl InChI=1/C14H19ClO3/c1-9(2)8-17-14(16)11(4)18-13-6-5-12(15)7-10(13)3/h5-7,9,11H,8H2,1-4H3/t11-/m1/s1 mecoprop-P-potassium InChI=1/C10H11ClO3.K/c1-6-5-8(11)3-4-9(6)14-7(2)10(12)13;/h3-5,7H,1-2H3,(H,12,13);/q;+1/p-1/t7-;/m1./s1/fC10H10ClO3.K/q-1;m			
	NOTE It should be stated which ester or salt is present, for example mecoprop-P-isobutyl [101012-85-5] or mecoprop-P-potassium [66423-05-0]. NOTE Il convient de préciser quel est l'ester ou le sel présent, par exemple mécoprop-P-isobutyle [101012-85-5] ou mécoprop-P-potassium [66423-05-0].			
E mefenacet **F** méfénacet (m)	2-(1,3-benzothiazol-2-yloxy)-*N*-methylacetanilide *N*-méthyl-(benzothiazol-2-yl)oxy-acétanilide 2-(2-benzothiazolyloxy)-*N*-methyl-*N*-phenylacetamide	 $C_{16}H_{14}N_2O_2S$	73250-68-7	**H**
	InChI=1/C16H14N2O2S/c1-18(12-7-3-2-4-8-12)15(19)11-20-16-17-13-9-5-6-10-14(13)21-16/h2-10H,11H2,1H3			

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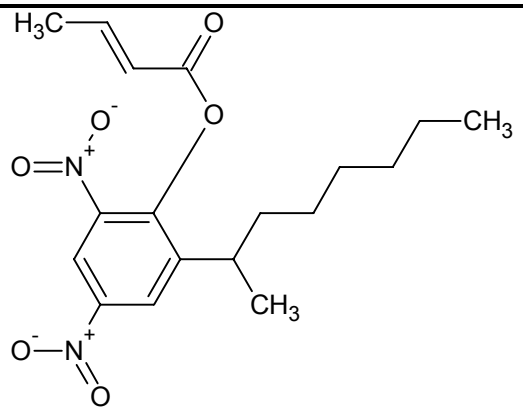
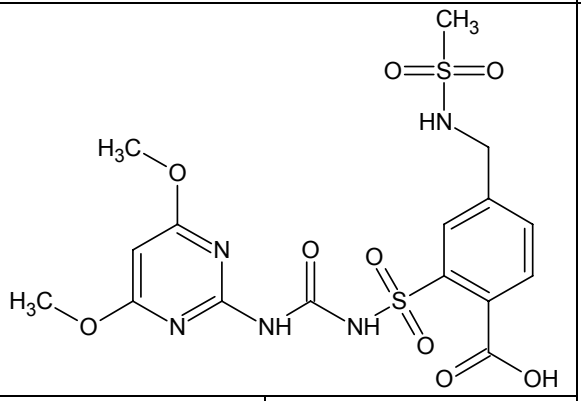
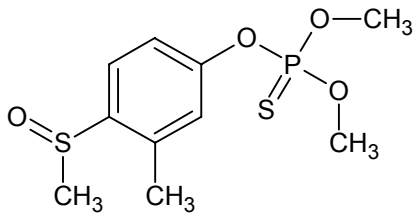
E meptyldinocap **F** méptyldinocap (m)	(*RS*)-2-(1-methylheptyl)-4,6-dinitrophenyl crotonate crotonate de (*RS*)-2-(oct-2-yl)-4,6-dinitrophényle 2-(1-methylheptyl)-4,6-dinitrophenyl (2*E*)-2-butenoate			**F**
		$C_{18}H_{24}N_2O_6$	131-72-6	
	InChI=1/C18H24N2O6/c1-4-6-7-8-10-13(3)15-11-14(19(22)23)12-16(20(24)25)18(15)26-17(21)9-5-2/h5,9,11-13H,4,6-8,10H2,1-3H3/b9-5+/t13-/s3			
E mesosulfuron **F** mésosulfuron (m)	2-[(4,6-dimethoxypyrimidin-2-ylcarbamoyl)sulfamoyl]-α-(methanesulfonamido)-*p*-toluic acid acide 2-[(4,6-diméthoxypyrimidin-2-yl-carbamoyl)sulfamoyl]-4-((méthylsulfamoyl)méthyl)-benzoïque 2-[[[[(4,6-dimethoxy-2-pyrimidinyl)=amino]carbonyl]amino]sulfonyl]-4-[[(methylsulfonyl)amino]methyl]=benzoic acid			**H**
		$C_{16}H_{19}N_5O_9S_2$	400852-66-6	
	InChI=1/C16H19N5O9S2/c1-29-12-7-13(30-2)19-15(18-12)20-16(24)21-32(27,28)11-6-9(8-17-31(3,25)26)4-5-10(11)14(22)23/h4-7,17H,8H2,1-3H3,(H,22,23)(H2,18,19,20,21,24)/f/h20-22H mesosulfuron-methyl InChI=1/C17H21N5O9S2/c1-29-13-8-14(30-2)20-16(19-13)21-17(24)22-33(27,28)12-7-10(9-18-32(4,25)26)5-6-11(12)15(23)31-3/h5-8,18H,9H2,1-4H3,(H2,19,20,21,22,24)/f/h21-22H			
	NOTE It should be stated which ester or salt is present, for example mesosulfuron-methyl [208465-21-8]. NOTE Il convient de préciser quel est l'ester ou le sel présent, par exemple mésosulfuron-méthyle [208465-21-8].			
E mesulfenfos **F** mésulfenfos (m)	*O,O*-dimethyl *O*-4-methylsulfinyl-*m*-tolyl phosphorothioate phosphorothioate de *O,O*-diméthyle et de *O*-(3-méthyl-4-sulfinylphényle) *O,O*-dimethyl *O*-[3-methyl-4-(methylsulfinyl)phenyl] phosphorothioate			**I**
		$C_{10}H_{15}O_4PS_2$	3761-41-9	
	InChI=1/C10H15O4PS2/c1-8-7-9(5-6-10(8)17(4)11)14-15(16,12-2)13-3/h5-7H,1-4H3			

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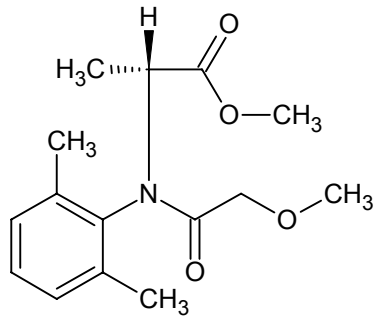
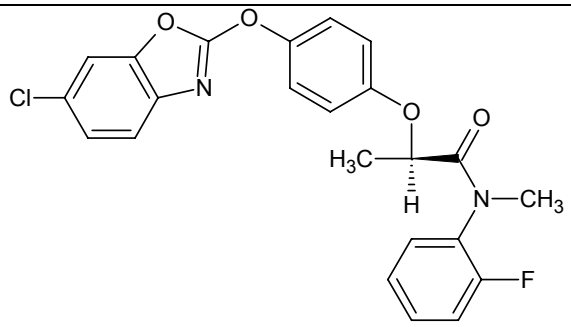
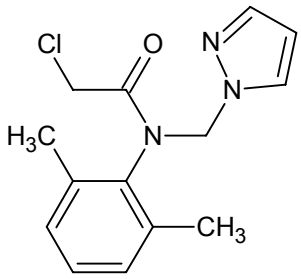
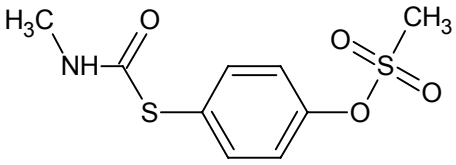
E metalaxyl-M **F** métalaxyl-M (m)	methyl *N*-(methoxyacetyl)-*N*-(2,6-xylyl)-D-alaninate D-(*N*-méthoxyacétyl-*N*-2,6-xylyl)alaninate de méthyle methyl *N*-(2,6-dimethylphenyl)-*N*-(methoxyacetyl)-D-alaninate			**F**
		$C_{15}H_{21}NO_4$	70630-17-0	
	InChI=1/C15H21NO4/c1-10-7-6-8-11(2)14(10)16(13(17)9-19-4)12(3)15(18)20-5/h6-8,12H,9H2,1-5H3/t12-/m1/s1			
E metamifop **F** métamifop (m)	(*R*)-2-[4-(6-chloro-1,3-benzoxazol-2-yloxy)phenoxy]-2′-fluoro-*N*-methylpropionanilide (*R*)-2-[4-(6-chloro-1,3-benzoxazol-2-yloxy)phénoxy]-*N*-(2-fluorophényl)-*N*-méthylpropanamide (2*R*)-2-[4-[(6-chloro-2-benzoxazolyl)oxy]phenoxy]-*N*-(2-fluorophenyl)-*N*-methylpropanamide			**H**
		$C_{23}H_{18}ClFN_2O_4$	256412-89-2	
	InChI=1/C23H18ClFN2O4/c1-14(22(28)27(2)20-6-4-3-5-18(20)25)29-16-8-10-17(11-9-16)30-23-26-19-12-7-15(24)13-21(19)31-23/h3-14H,1-2H3/t14-/m1/s1			
E metazachlor **F** métazachlore (m)	2-chloro-*N*-(pyrazol-1-ylmethyl)acet-2′,6′-xylidide 2-chloro-2′,6′-diméthyl-*N*-[(pyrazol-1-yl)méthyl]acétanilide 2-chloro-*N*-(2,6-dimethylphenyl)-*N*-(1*H*-pyrazol-1-ylmethyl)acetamide			**H**
		$C_{14}H_{16}ClN_3O$	67129-08-2	
	InChI=1/C14H16ClN3O/c1-11-5-3-6-12(2)14(11)18(13(19)9-15)10-17-8-4-7-16-17/h3-8H,9-10H2,1-2H3			
E methasulfocarb **F** méthasulfocarb (m)	*S*-4-(mesyloxy)phenyl methyl(thiocarbamate) méthylthiocarbamate de *S*-4-(mésyloxy)phényle *S*-[4-[(methylsulfonyl)oxy]phenyl] methylcarbamothioate			**F** **P**
		$C_9H_{11}NO_4S_2$	66952-49-6	
	InChI=1/C9H11NO4S2/c1-10-9(11)15-8-5-3-7(4-6-8)14-16(2,12)13/h3-6H,1-2H3,(H,10,11)/f/h10H			

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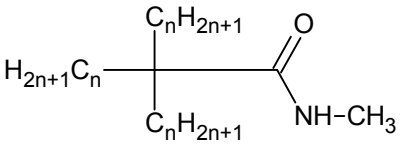
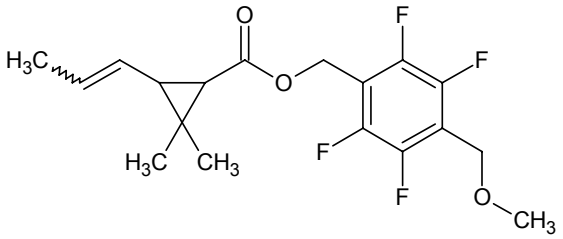
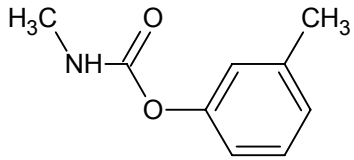
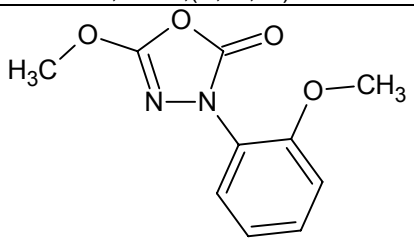
	IUPAC International Chemical Identifier (InChI™)			
E methylneodecanamide **F** méthylnéodécanamide (m)	2,2,2-trialkyl-*N*-methylacetamide *N*-méthylnéodécanamide *N*-methylneodecanamide			**RE**
		n/a	105726-67-8	
	n/a			
E metofluthrin **F** métofluthrine (f)	2,3,5,6-tetrafluoro-4-(methoxymethyl)benzyl (*EZ*)-(1*RS*,3*RS*;1*RS*,3*SR*)-2,2-dimethyl-3-prop-1-enylcyclopropanecarboxylate Alternative Rothamsted-style stereodescriptors: 2,3,5,6-tetrafluoro-4-(methoxymethyl)benzyl (*EZ*)-(1*RS*)-*cis-trans*-2,2-dimethyl-3-prop-1-enylcyclopropanecarboxylate (*EZ*)-(1*RS*,3*RS;1RS,3SR)*-2,2-diméthyl-3-(prop-1-ènyl)cyclopropane-carboxylate de (2,3,5,6-tétrafluoro-4-(méthoxyméthyl)benzyle [2,3,5,6-tetrafluoro-4-(methoxymethyl)phenyl]methyl 2,2-dimethyl-3-(1-propenyl)cyclopropanecarboxylate			**I**
		$C_{18}H_{20}F_4O_3$	240494-70-6	
	InChI=1/C18H20F4O3/c1-5-6-11-12(18(11,2)3)17(23)25-8-10-15(21)13(19)9(7-24-4)14(20)16(10)22/h5-6,11-12H,7-8H2,1-4H3/b6-5?/t11?,12?			
E metolcarb **F** métolcarb (m)	*m*-tolyl methylcarbamate méthylcarbamate de *m*-tolyle 3-methylphenyl methylcarbamate			**A** **I**
		$C_9H_{11}NO_2$	1129-41-5	
	InChI=1/C9H11NO2/c1-7-4-3-5-8(6-7)12-9(11)10-2/h3-6H,1-2H3,(H,10,11)/f/h10H			
E metoxadiazone **F** métoxadiazone (f)	5-methoxy-3-(2-methoxyphenyl)-1,3,4-oxadiazol-2(3*H*)-one 5-méthoxy-3-(2-méthoxyphényl)-1,3,4-(3*H*)-oxadiazol-2-one 5-methoxy-3-(2-methoxyphenyl)-1,3,4-oxadiazol-2(3*H*)-one			**I**
		$C_{10}H_{10}N_2O_4$	60589-06-2	
	InChI=1/C10H10N2O4/c1-14-8-6-4-3-5-7(8)12-10(13)16-9(11-12)15-2/h3-6H,1-2H3			

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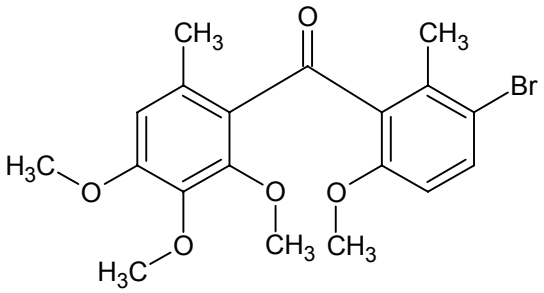
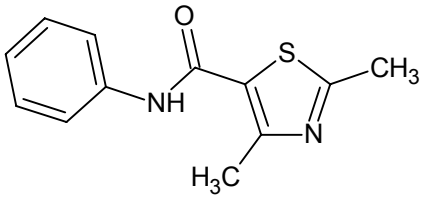
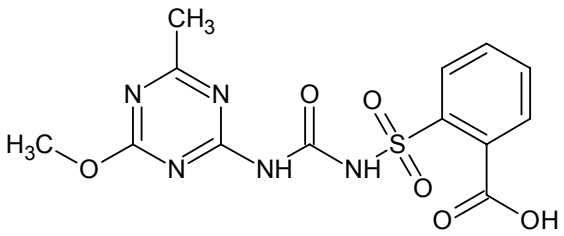
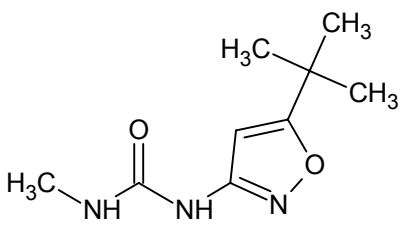
E metrafenone **F** métrafénone (f)	3′-bromo-2,3,4,6′-tetramethoxy-2′,6-dimethylbenzophenone (3-bromo-6-méthoxy-2-méthylphényl)-(2,3,4-triméthoxy-6-méthylphényl)méthanone (3-bromo-6-methoxy-2-methylphenyl)(2,3,4-trimethoxy-6-methylphenyl)methanone			**F**
		$C_{19}H_{21}BrO_5$	220899-03-6	
	InChI=1/C19H21BrO5/c1-10-9-14(23-4)18(24-5)19(25-6)15(10)17(21)16-11(2)12(20)7-8-13(16)22-3/h7-9H,1-6H3			
E metsulfovax **F** metsulfovax (m)	2,4-dimethyl-1,3-thiazole-5-carboxanilide 2,4-diméthyl-1,3-thiazol-5-ylcarboxanilide 2,4-dimethyl-*N*-phenyl-5-thiazolecarboxamide			**F**
		$C_{12}H_{12}N_2OS$	21452-18-6	
	InChI=1/C12H12N2OS/c1-8-11(16-9(2)13-8)12(15)14-10-6-4-3-5-7-10/h3-7H,1-2H3,(H,14,15)/f/h14H			
E metsulfuron **F** metsulfuron (m)	2-(4-methoxy-6-methyl-1,3,5-triazin-2-ylcarbamoylsulfamoyl)benzoic acid acide 2-(4-méthoxy-6-méthyl-1,3,5-triazin-2-yl)-uréidosulfonylbenzoïque 2-[[[[(4-methoxy-6-methyl-1,3,5-triazin-2-yl)amino]carbonyl]=amino]sulfonyl]benzoic acid			**H**
		$C_{13}H_{13}N_5O_6S$	79510-48-8	
	InChI=1/C13H13N5O6S/c1-7-14-11(17-13(15-7)24-2)16-12(21)18-25(22,23)9-6-4-3-5-8(9)10(19)20/h3-6H,1-2H3,(H,19,20)(H2,14,15,16,17,18,21)/f/h16,18-19H metsulfuron-methyl InChI=1/C14H15N5O6S/c1-8-15-12(18-14(16-8)25-3)17-13(21)19-26(22,23)10-7-5-4-6-9(10)11(20)24-2/h4-7H,1-3H3,(H2,15,16,17,18,19,21)/f/h17,19H			
	NOTE It should be stated which ester or salt is present, for example metsulfuron-methyl [74223-64-6]. NOTE Il convient de préciser quel est l'ester ou le sel présent, par exemple metsulfuron-méthyle [74223-64-6].			
E monisouron **F** monisuron (m)	1-(5-*tert*-butyl-1,2-oxazol-3-yl)-3-methylurea *N*-(5-*tert*-butyl-1,2-oxazol-3-yl)-*N*′-méthylurée ou 1-(5-*tert*-butyl-1,2-oxazol-3-yl)-3-méthylurée *N*-[5-(1,1-dimethylethyl)-3-isoxazolyl]-*N*′-methylurea			**H**
		$C_9H_{15}N_3O_2$	55807-46-0	
	InChI=1/C9H15N3O2/c1-9(2,3)6-5-7(12-14-6)11-8(13)10-4/h5H,1-4H3,(H2,10,11,12,13)/f/h10-11H			

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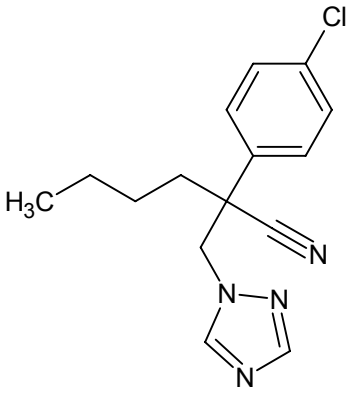
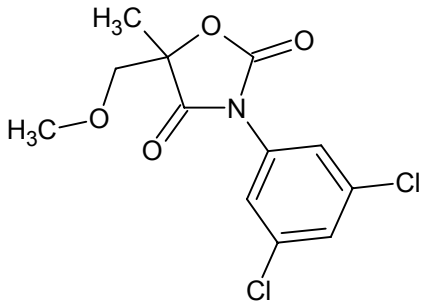
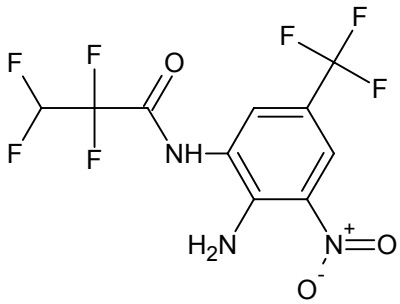
		Molecular formula Formule brute	CAS Registry Number® Numéro d'enregistrement CAS®	Application
	IUPAC International Chemical Identifier (InChI™)			
E myclobutanil **F** myclobutanil (m)	(*RS*)-2-(4-chlorophenyl)-2-(1*H*-1,2,4-triazol-1-ylmethyl)hexanenitrile (*RS*)-2-(4-chlorophényl)-2-(((1*H*)-1,2,4-triazol-1-yl)méthyl)hexanenitrile α-butyl-α-(4-chlorophenyl)-1*H*-1,2,4-triazole-1-propanenitrile			**F**
		$C_{15}H_{17}ClN_4$	88671-89-0	
	InChI=1/C15H17ClN4/c1-2-3-8-15(9-17,10-20-12-18-11-19-20)13-4-6-14(16)7-5-13/h4-7,11-12H,2-3,8,10H2,1H3/t15-/s3			
E myclozolin **F** myclozoline (f)	(*RS*)-3-(3,5-dichlorophenyl)-5-methoxymethyl-5-methyl-1,3-oxazolidine-2,4-dione (*RS*)-3-(3,5-dichlorophényl)-5-méthoxyméthyl-5-méthyl-1,3-oxazolidine-2,4-dione 3-(3,5-dichlorophenyl)-5-(methoxymethyl)-5-methyl-2,4-oxazolidinedione			**F**
		$C_{12}H_{11}Cl_2NO_4$	54864-61-8	
	InChI=1/C12H11Cl2NO4/c1-12(6-18-2)10(16)15(11(17)19-12)9-4-7(13)3-8(14)5-9/h3-5H,6H2,1-2H3/t12-/s3			
E nifluridide **F** nifluridide (m)	6′-amino-α,α,α,2,2,3,3-heptafluoro-5′-nitropropion-*m*-toluidide 2′-amino-3′-nitro-5′-trifluorométhyl-2,2,3,3-tétrafluoropropionanilide *N*-[2-amino-3-nitro-5-(trifluoromethyl)phenyl]-2,2,3,3-tetrafluoropropanamide			**A** **I**
		$C_{10}H_6F_7N_3O_3$	61444-62-0	
	InChI=1/C10H6F7N3O3/c11-7(12)9(13,14)8(21)19-4-1-3(10(15,16)17)2-5(6(4)18)20(22)23/h1-2,7H,18H2,(H,19,21)/f/h19H			

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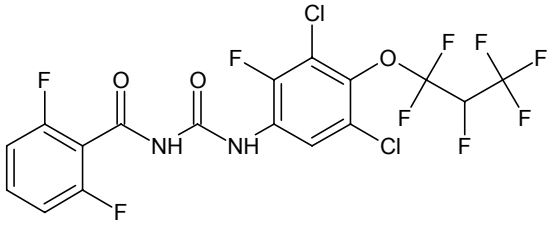
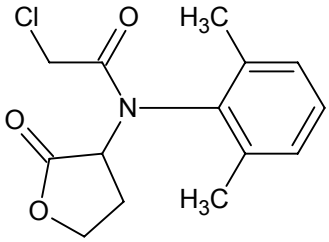
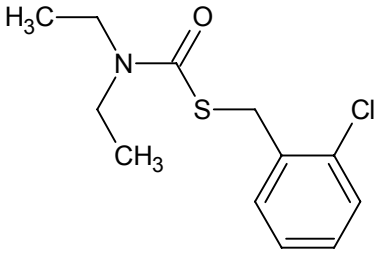
E noviflumuron **F** noviflumuron (m)	1-{3,5-dichloro-2-fluoro-4-[(*RS*)-1,1,2,3,3,3-hexafluoropropoxy]=phenyl}-3-(2,6-difluorobenzoyl)urea (*RS*)-*N*-[3,5-dichloro-2-fluoro-4-(1,1,2,3,3,3-hexafluoropropoxy)=phényl]-*N*′-(2,6-difluorobenzoyl)urée ou (*RS*)-1-[3,5-dichloro-2-fluoro-4-(1,1,2,3,3,3-hexafluoropropoxy)=phényl]-3-(2,6-difluorobenzoyl)urée *N*-[[[3,5-dichloro-2-fluoro-4-(1,1,2,3,3,3-hexafluoropropoxy)=phenyl]amino]carbonyl]-2,6-difluorobenzamide			**IGR**
		$C_{17}H_7Cl_2F_9N_2O_3$	121451-02-3	
	InChI=1/C17H7Cl2F9N2O3/c18-5-4-8(29-15(32)30-13(31)9-6(20)2-1-3-7(9)21)11(22)10(19)12(5)33-17(27,28)14(23)16(24,25)26/h1-4,14H,(H2,29,30,31,32)/t14-/s3/f/h29-30H			
E ofurace **F** ofurace (m)	(*RS*)-α-(2-chloro-*N*-2,6-xylylacetamido)-γ-butyrolactone *N*-(*RS*)-(γ-butyrolacton-3-yl)-2-chloro-2′,6′-diméthylacétanilide 2-chloro-*N*-(2,6-dimethylphenyl)-*N*-(tetrahydro-2-oxo-3-furanyl)acetamide			**F**
		$C_{14}H_{16}ClNO_3$	58810-48-3	
	InChI=1/C14H16ClNO3/c1-9-4-3-5-10(2)13(9)16(12(17)8-15)11-6-7-19-14(11)18/h3-5,11H,6-8H2,1-2H3/t11-/s3			
E orbencarb **F** orbencarb (m)	*S*-2-chlorobenzyl diethyl(thiocarbamate) diéthylthiocarbamate de *S*-2-chlorobenzyle *S*-[(2-chlorophenyl)methyl] diethylcarbamothioate			**H**
		$C_{12}H_{16}ClNOS$	34622-58-7	
	InChI=1/C12H16ClNOS/c1-3-14(4-2)12(15)16-9-10-7-5-6-8-11(10)13/h5-8H,3-4,9H2,1-2H3			

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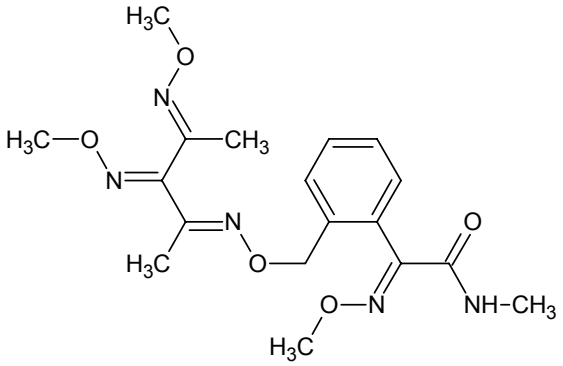
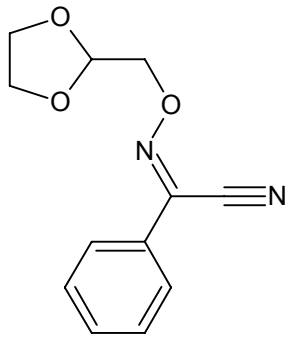
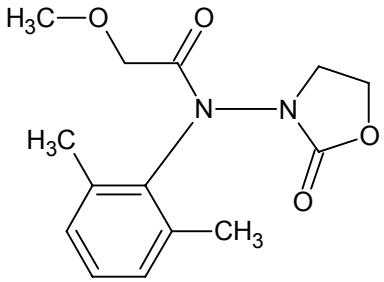
E orysastrobin **F** orysastrobine (f)	(2*E*)-2-(methoxyimino)-2-{2-[(3*E*,5*E*,6*E*)-5-(methoxyimino)-4,6-dimethyl-2,8-dioxa-3,7-diazanona-3,6-dien-1-yl]phenyl}-*N*-methylacetamide (2*E*)-2-(méthoxyimino)-2-{2-[(3*E*,5*E*,6*E*)-5-(méthoxyimino)-4,6-diméthyl-2,8-dioxa-3,7-diazanona-3,6-dién-1-yl]phényl}-*N*-méthylacétamide (α*E*)-α-(methoxyimino)-2-[(3*E*,5*E*,6*E*)-5-(methoxyimino)-4,6-dimethyl-2,8-dioxa-3,7-diaza-3,6-nonadienyl]-*N*-methylbenzeneacetamide			**F**
		$C_{18}H_{25}N_5O_5$	248593-16-0	
	InChI=1/C18H25N5O5/c1-12(20-25-4)16(22-26-5)13(2)21-28-11-14-9-7-8-10-15(14)17(23-27-6)18(24)19-3/h7-10H,11H2,1-6H3,(H,19,24)/b20-12+,21-13+,22-16+,23-17+/f/h19H			
E oxabetrinil **F** oxabétrinil (m)	(*Z*)-1,3-dioxolan-2-ylmethoxyimino(phenyl)acetonitrile (*Z*)-(1,3-dioxolan-2-yl(méthoxy)imino)-phénylacétonitrile (α*Z*)-α-[(1,3-dioxolan-2-yl)methoxyimino]benzeneacetonitrile			**S**
		$C_{12}H_{12}N_2O_3$	94593-79-0	
	InChI=1/C12H12N2O3/c13-8-11(10-4-2-1-3-5-10)14-17-9-12-15-6-7-16-12/h1-5,12H,6-7,9H2/b14-11+			
E oxadixyl **F** oxadixyl (m)	2-methoxy-*N*-(2-oxo-1,3-oxazolidin-3-yl)acet-2′,6′-xylidide 2-méthoxy-*N*-(2-oxo-1,3-oxazolidin-3-yl)-2′,6′-diméthylacétanilide *N*-(2,6-dimethylphenyl)-2-methoxy-*N*-(2-oxo-3-oxazolidinyl)acetamide			**F**
		$C_{14}H_{18}N_2O_4$	77732-09-3	
	InChI=1/C14H18N2O4/c1-10-5-4-6-11(2)13(10)16(12(17)9-19-3)15-7-8-20-14(15)18/h4-6H,7-9H2,1-3H3			

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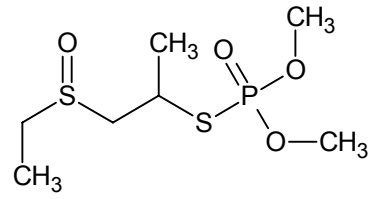
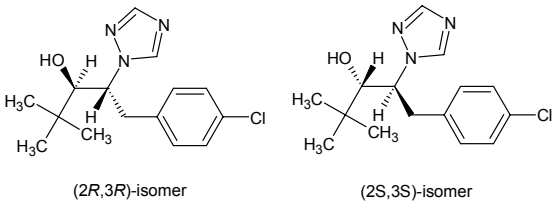
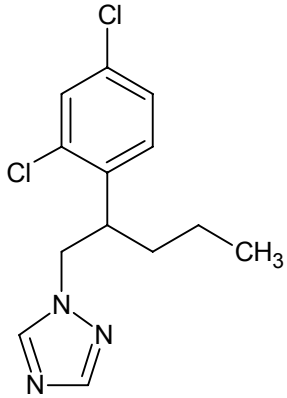
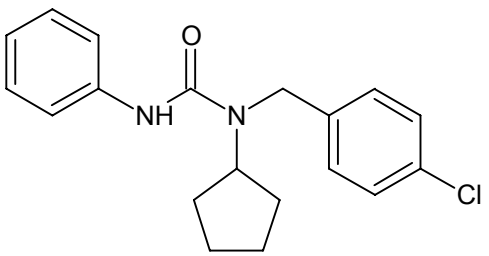
E oxydeprofos **F** oxydéprofos (m)	*S*-(*RS*)-2-ethylsulfinyl-1-methylethyl *O,O*-dimethyl phosphorothioate phosphorothioate de *S*-(*RS*)-2-éthylsulfinyl-1-méthyléthyle et de *O,O*-diméthyle *S*-[2-(ethylsulfinyl)-1-methylethyl] *O,O*-dimethyl phosphorothioate	 $C_7H_{17}O_4PS_2$	2674-91-1	**A** **I**
	InChI=1/C7H17O4PS2/c1-5-14(9)6-7(2)13-12(8,10-3)11-4/h7H,5-6H2,1-4H3			
E paclobutrazol **F** paclobutrazol (m)	(2*RS*,3*RS*)-1-(4-chlorophenyl)-4,4-dimethyl-2-(1*H*-1,2,4-triazol-1-yl)pentan-3-ol (2*RS*,3*RS*)-1-(4-chlorophényl)-4,4-diméthyl-2-(1*H*-1,2,4-triazol-1-yl)pentan-3-ol (α*R*,β*R*)-*rel*-β-[(4-chlorophenyl)=methyl]-α-(1,1-dimethylethyl)-1*H*-1,2,4-triazole-1-ethanol	 (2*R*,3*R*)-isomer (2S,3S)-isomer $C_{15}H_{20}ClN_3O$	76738-62-0	**P**
	InChI=1/C15H20ClN3O/c1-15(2,3)14(20)13(19-10-17-9-18-19)8-11-4-6-12(16)7-5-11/h4-7,9-10,13-14,20H,8H2,1-3H3/t13-,14+/s3			
E penconazole **F** penconazole (m)	(*RS*)-1-[2-(2,4-dichlorophenyl)pentyl]-1*H*-1,2,4-triazole (*RS*)-1-[2-(2,4-dichlorophényl)pentyl]-1*H*-1,2,4-triazole 1-[2-(2,4-dichlorophenyl)pentyl]-1*H*-1,2,4-triazole	 $C_{13}H_{15}Cl_2N_3$	66246-88-6	**F**
	InChI=1/C13H15Cl2N3/c1-2-3-10(7-18-9-16-8-17-18)12-5-4-11(14)6-13(12)15/h4-6,8-10H,2-3,7H2,1H3/t10-/s3			
E pencycuron **F** pencycuron (m)	1-(4-chlorobenzyl)-1-cyclopentyl-3-phenylurea *N*-(4-chlorobenzyl)-*N*-cyclopentyl-*N′*-phénylurée ou 1-(4-chlorobenzyl)-1-cyclopentyl-3-phénylurée *N*-[(4-chlorophenyl)methyl]-*N*-cyclopentyl-*N′*-phenylurea	 $C_{19}H_{21}ClN_2O$	66063-05-6	**F**
	InChI=1/C19H21ClN2O/c20-16-12-10-15(11-13-16)14-22(18-8-4-5-9-18)19(23)21-17-6-2-1-3-7-17/h1-3,6-7,10-13,18H,4-5,8-9,14H2,(H,21,23)/f/h21H			

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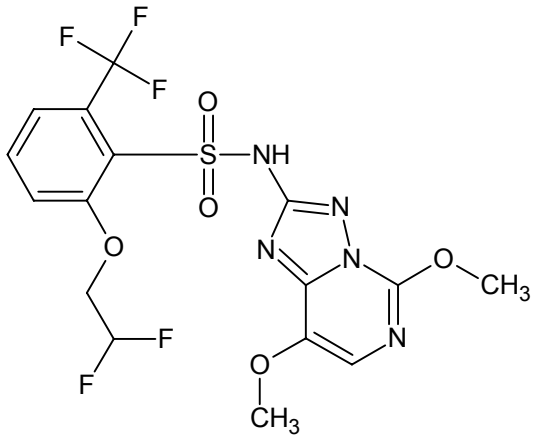
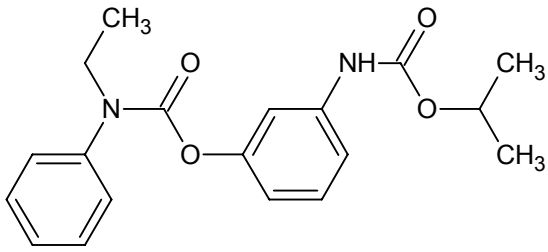
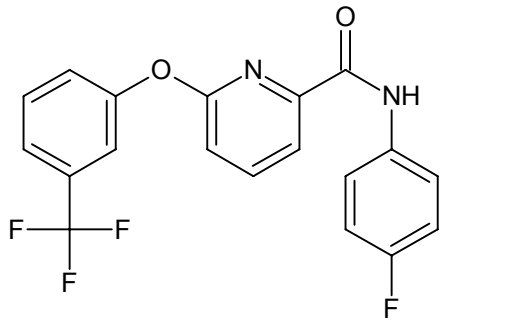
E penoxsulam **F** pénoxsulame (m)	3-(2,2-difluoroethoxy)-*N*-(5,8-dimethoxy[1,2,4]triazolo[1,5-*c*]pyrimidin-2-yl)-α,α,α-trifluorotoluene-2-sulfonamide 2-(2,2-difluoroéthoxy)-*N*-(5,8-diméthoxy-1,2,4-triazolo[1,5-*c*]pyrimidin-3-yl)-6-trifluorométhylbenzène-sulfamide 2-(2,2-difluoroethoxy)-*N*-(5,8-dimethoxy[1,2,4]triazolo[1,5-*c*]pyrimidin-2-yl)-6-(trifluoromethyl)benzenesulfonamide	 $C_{16}H_{14}F_5N_5O_5S$	219714-96-2	**H**
	InChI=1/C16H14F5N5O5S/c1-29-10-6-22-15(30-2)26-13(10)23-14(24-26)25-32(27,28)12-8(16(19,20)21)4-3-5-9(12)31-7-11(17)18/h3-6,11H,7H2,1-2H3,(H,24,25)/f/h25H			
E phenisopham **F** phénisophame (m)	3-(isopropoxycarbonylamino)phenyl *N*-ethylcarbanilate *N*-éthylcarbanilate de 3-(isopropoxycarbonylamino)phényle 3-[[(1-methylethoxy)carbonyl]amino]=phenyl ethylphenylcarbamate	 $C_{19}H_{22}N_2O_4$	57375-63-0	**H**
	InChI=1/C19H22N2O4/c1-4-21(16-10-6-5-7-11-16)19(23)25-17-12-8-9-15(13-17)20-18(22)24-14(2)3/h5-14H,4H2,1-3H3,(H,20,22)/f/h20H			
E picolinafen **F** picolinafène (m)	4′-fluoro-6-(α,α,α-trifluoro-*m*-tolyloxy)pyridine-2-carboxanilide *N*-(4-fluorophényl)-6-(3-trifluorométhyl)phénoxypyridine-2-carboxamide *N*-(4-fluorophenyl)-6-[3-(trifluoromethyl)phenoxy]-2-pyridinecarboxamide	 $C_{19}H_{12}F_4N_2O_2$	137641-05-5	**H**
	InChI=1/C19H12F4N2O2/c20-13-7-9-14(10-8-13)24-18(26)16-5-2-6-17(25-16)27-15-4-1-3-12(11-15)19(21,22)23/h1-11H,(H,24,26)/f/h24H			

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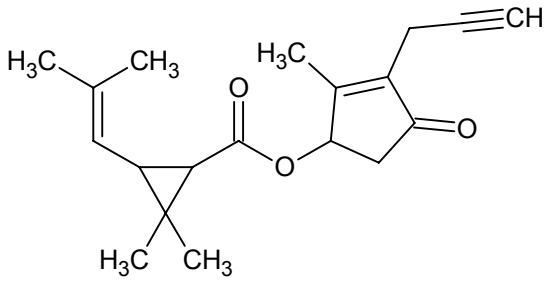
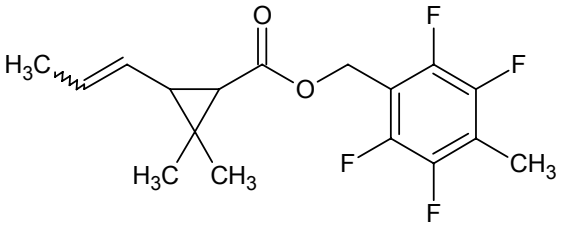
E prallethrin **F** pralléthrine (f)	(*RS*)-2-methyl-4-oxo-3-prop-2-ynylcyclopent-2-enyl (1*RS*,3*RS*;1*RS*,3*SR*)-2,2-dimethyl-3-(2-methylprop-1-enyl)cyclopropanecarboxylate Alternative Rothamsted-style stereodescriptors: (*RS*)-2-methyl-4-oxo-3-prop-2-ynylcyclopent-2-enyl (1*RS*)-*cis-trans*-2,2-dimethyl-3-(2-methylprop-1-enyl)cyclopropanecarboxylate			**I**
	(1*RS*,3*RS*;1*RS*,3*SR*)-2,2-diméthyl-3-(2-méthylprop-1-ènyl)-cyclopropanecarboxylate de (*RS*)-(2-méthyl-4-oxo-3-prop-2-ynyl)cyclopent-2-èn-1-yle			
	2-methyl-4-oxo-3-(2-propynyl)-2-cyclopenten-1-yl 2,2-dimethyl-3-(2-methyl-1-propenyl)cyclopropanecarboxylate	$C_{19}H_{24}O_3$	23031-36-9	
	InChI=1/C19H24O3/c1-7-8-13-12(4)16(10-15(13)20)22-18(21)17-14(9-11(2)3)19(17,5)6/h1,9,14,16-17H,8,10H2,2-6H3/t14?,16?,17?			
	NOTE The name “prallethrin” is not acceptable for use in France because it does not reflect the chemical name. NOTE Le nom «pralléthrine» n'est pas acceptable pour l'emploi en France, parce qu'il est sans lien avec la formule chimique.			
E profluthrin **F** profluthrine (f)	2,3,5,6-tetrafluoro-4-methylbenzyl (*EZ*)-(1*RS*,3*RS*;1*RS*,3*SR*)-2,2-dimethyl-3-prop-1-enylcyclopropanecarboxylate Alternative Rothamsted-style stereodescriptors: 2,3,5,6-tetrafluoro-4-methylbenzyl (*EZ*)-(1*RS*)-*cis-trans*-2,2-dimethyl-3-prop-1-enylcyclopropanecarboxylate			**I**
	(*EZ*)-(1*RS*,3*RS*;1*RS*,3*SR*)-2,2-diméthyl-3-(prop-1-ènyl)cyclopropanecarboxylate de 2,3,5,6-tétrafluoro-4-méthylbenzyle			
	(2,3,5,6-tetrafluoro-4-methylphenyl)methyl 2,2-dimethyl-3-(1-propenyl)cyclopropanecarboxylate	$C_{17}H_{18}F_4O_2$	223419-20-3	
	InChI=1/C17H18F4O2/c1-5-6-10-11(17(10,3)4)16(22)23-7-9-14(20)12(18)8(2)13(19)15(9)21/h5-6,10-11H,7H2,1-4H3/b6-5?/t10?,11?			

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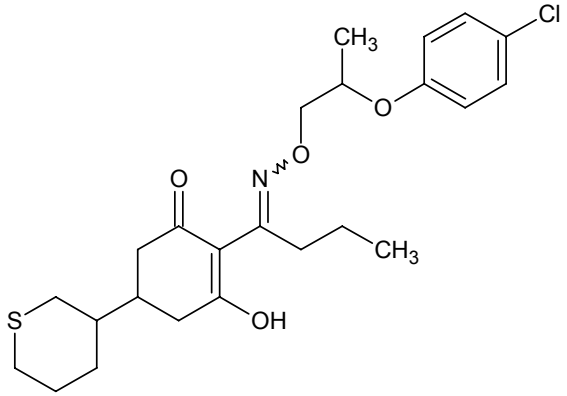
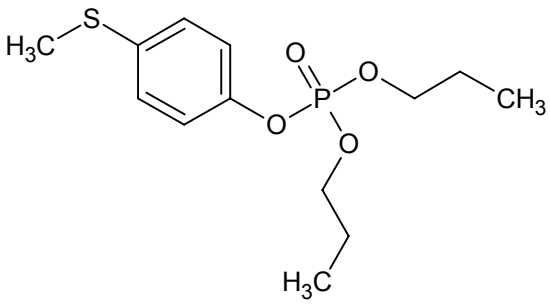
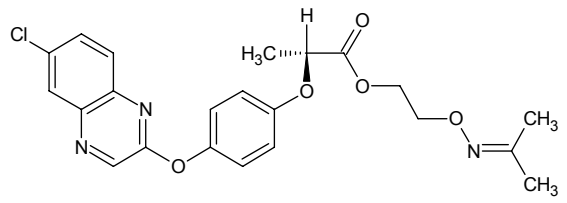
E profoxydim **F** profoxydime (m)	(5*RS*)-2-{(*EZ*)-1-[(2*RS*)-2-(4-chlorophenoxy)propoxyimino]butyl}-3-hydroxy-5-[(3*RS*)-thian-3-yl]cyclohex-2-en-1-one (5*RS*)-2-{(*EZ*)-1-[(2*RS*)-2-(4-chlorophénoxy)propoxyimino]butyl}-3-hydroxy-5-[(3*RS*)-tétrahydro-(2*H*)thiopyran-3-yl]cyclohex-2-énone 2-[1-[[2-(4-chlorophenoxy)propoxy]imino]butyl]-3-hydroxy-5-(tetrahydro-2*H*-thiopyran-3-yl)-2-cyclohexen-1-one	 $C_{24}H_{32}ClNO_4S$	139001-49-3	**H**
	InChI=1/C24H32ClNO4S/c1-3-5-21(26-29-14-16(2)30-20-9-7-19(25)8-10-20)24-22(27)12-18(13-23(24)28)17-6-4-11-31-15-17/h7-10,16-18,27H,3-6,11-15H2,1-2H3/b26-21?/t16?,17?,18?			
E propaphos **F** propafos (m)	4-(methylthio)phenyl dipropyl phosphate phosphate de 4-(méthylthio)phényle et de dipropyle 4-(methylthio)phenyl dipropyl phosphate	 $C_{13}H_{21}O_4PS$	7292-16-2	**I**
	InChI=1/C13H21O4PS/c1-4-10-15-18(14,16-11-5-2)17-12-6-8-13(19-3)9-7-12/h6-9H,4-5,10-11H2,1-3H3			
E propaquizafop **F** propaquizafop (m)	2-isopropylideneamino-oxyethyl (*R*)-2-[4-(6-chloroquinoxalin-2-yloxy)phenoxy]propionate (*R*)-2-[4-(6-chloroquinoxalin-2-yloxy)phénoxy]propionate de 2-[(isopropylidène)amino]oxy]éthyle 2-[[(1-methylethylidene)amino]oxy]ethyl (2*R*)-2-[4-[(6-chloro-2-quinoxalinyl)oxy]phenoxy]propanoate	 $C_{22}H_{22}ClN_3O_5$	111479-05-1	**H**
	InChI=1/C22H22ClN3O5/c1-14(2)26-29-11-10-28-22(27)15(3)30-17-5-7-18(8-6-17)31-21-13-24-20-12-16(23)4-9-19(20)25-21/h4-9,12-13,15H,10-11H2,1-3H3/t15-/m1/s1			

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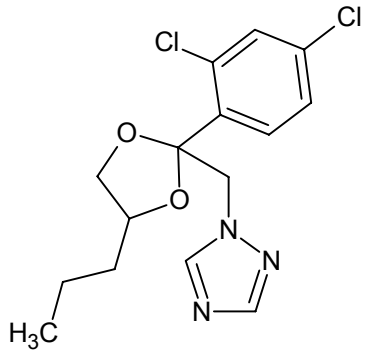
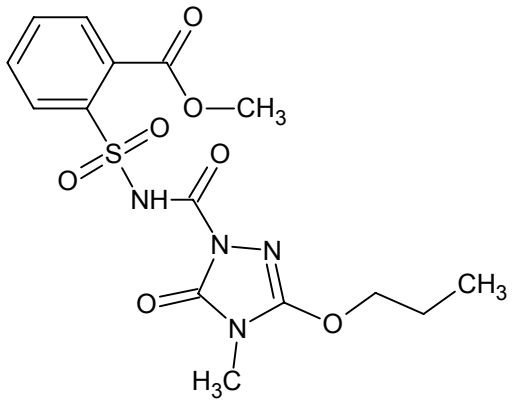
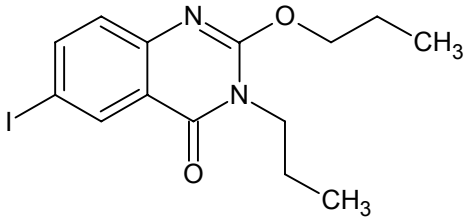
E propiconazole **F** propiconazole (m)	(2*RS*,4*RS*;2*RS*,4*SR*)-1-[2-(2,4-dichlorophenyl)-4-propyl-1,3-dioxolan-2-ylmethyl]-1*H*-1,2,4-triazole (2*RS*,4*RS*;2*RS*,4*SR*)-1-[2-(2,4-dichlorophényl)-4-propyl-1,3-dioxolan-2-yl]-méthyl-1*H*-1,2,4-triazole 1-[[2-(2,4-dichlorophenyl)-4-propyl-1,3-dioxolan-2-yl]methyl]-1*H*-1,2,4-triazole			**F**
		$C_{15}H_{17}Cl_2N_3O_2$	60207-90-1	
	InChI=1/C15H17Cl2N3O2/c1-2-3-12-7-21-15(22-12,8-20-10-18-9-19-20)13-5-4-11(16)6-14(13)17/h4-6,9-10,12H,2-3,7-8H2,1H3/t12?,15?			
E propoxycarbazone **F** propoxycarbazone (f)	methyl 2-[(4,5-dihydro-4-methyl-5-oxo-3-propoxy-1*H*-1,2,4-triazole-1-carboxamido)sulfonyl]benzoate 2-[(4-méthyl-5-oxo-3-propoxy)-(4,5-dihydro-1*H*-1,2,4-triazol-1-yl)]carboxamidosulfonyl]benzoate de méthyle methyl 2-[[[(4,5-dihydro-4-methyl-5-oxo-3-propoxy-1*H*-1,2,4-triazol-1-yl)carbonyl]amino]sulfonyl]benzoate			**H**
		$C_{15}H_{18}N_4O_7S$	145026-81-9	
	InChI=1/C15H18N4O7S/c1-4-9-26-14-16-19(15(22)18(14)2)13(21)17-27(23,24)11-8-6-5-7-10(11)12(20)25-3/h5-8H,4,9H2,1-3H3,(H,17,21)/f/h17H propoxycarbazone-sodium InChI=1/C15H18N4O7S.Na/c1-4-9-26-14-16-19(15(22)18(14)2)13(21)17-27(23,24)11-8-6-5-7-10(11)12(20)25-3;/h5-8H,4,9H2,1-3H3,(H,17,21);/q;+1/p-1/fC15H17N4O7S.Na/q-1;m/b17-13?;			
	NOTE It should be stated which salt is present, for example propoxycarbazone-sodium [181274-15-7]. NOTE Il convient de préciser quel est le sel présent, par exemple propoxycarbazone-sodium [181274-15-7].			
E proquinazid **F** proquinazide (m)	6-iodo-2-propoxy-3-propylquinazolin-4(3*H*)-one 6-iodo-2-propoxy-3-propyl-3*H*-quinazolin-4-one 6-iodo-2-propoxy-3-propyl-4(3*H*)-quinazolinone			**F**
		$C_{14}H_{17}IN_2O_2$	189278-12-4	
	InChI=1/C14H17IN2O2/c1-3-7-17-13(18)11-9-10(15)5-6-12(11)16-14(17)19-8-4-2/h5-6,9H,3-4,7-8H2,1-2H3			

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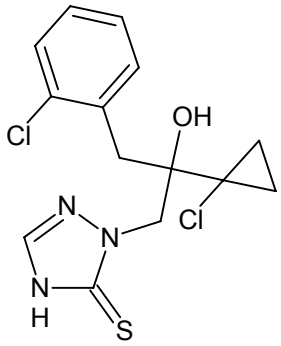
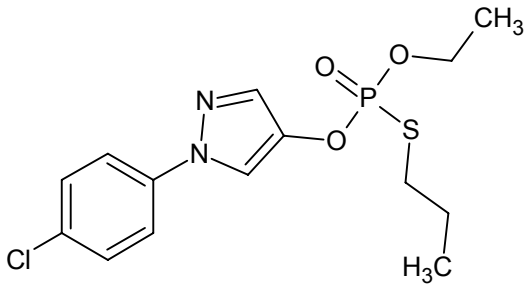
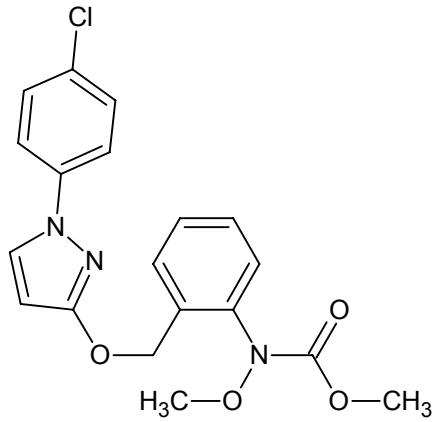
E prothioconazole **F** prothioconazole (m)	(*RS*)-2-[2-(1-chlorocyclopropyl)-3-(2-chlorophenyl)-2-hydroxypropyl]-2,4-dihydro-1,2,4-triazole-3-thione (*RS*)-2-[2-(1-chlorocyclopropyl)-3-(2-chlorophényl)-2-hydroxypropyl]-2,4-dihydro-1,2,4-triazole-3-thione 2-[2-(1-chlorocyclopropyl)-3-(2-chlorophenyl)-2-hydroxypropyl]-1,2-dihydro-3*H*-1,2,4-triazole-3-thione			**F**
		$C_{14}H_{15}Cl_2N_3OS$	178928-70-6	
	InChI=1/C14H15Cl2N3OS/c15-11-4-2-1-3-10(11)7-14(20,13(16)5-6-13)8-19-12(21)17-9-18-19/h1-4,9,20H,5-8H2,(H,17,18,21)/t14-/s3/f/h17H			
E pyraclofos **F** pyraclofos (m)	*O*-1-(4-chlorophenyl)pyrazol-4-yl *O*-ethyl *S*-propyl phosphorothioate phosphorothioate de *O*-1-(4-chlorophényl)pyrazol-4-yle de *O*-éthyle et de *S*-propyle *O*-[1-(4-chlorophenyl)-1*H*-pyrazol-4-yl] *O*-ethyl *S*-propyl phosphorothioate			**I**
		$C_{14}H_{18}ClN_2O_3PS$	77458-01-6	
	InChI=1/C14H18ClN2O3PS/c1-3-9-22-21(18,19-4-2)20-14-10-16-17(11-14)13-7-5-12(15)6-8-13/h5-8,10-11H,3-4,9H2,1-2H3/t21-/s3			
E pyraclostrobin **F** pyraclostrobine (f)	methyl {2-[1-(4-chlorophenyl)pyrazol-3-yloxymethyl]phenyl}=(methoxy)carbamate *N*-{2-[1-(4-chlorophényl)pyrazol-3-yloxyméthyl]phényl}-*N*-(méthoxy)carbamate de méthyle methyl [2-[[[1-(4-chlorophenyl)-1*H*-pyrazol-3-yl]oxy]methyl]phenyl]=methoxycarbamate			**F**
		$C_{19}H_{18}ClN_3O_4$	175013-18-0	
	InChI=1/C19H18ClN3O4/c1-25-19(24)23(26-2)17-6-4-3-5-14(17)13-27-18-11-12-22(21-18)16-9-7-15(20)8-10-16/h3-12H,13H2,1-2H3			

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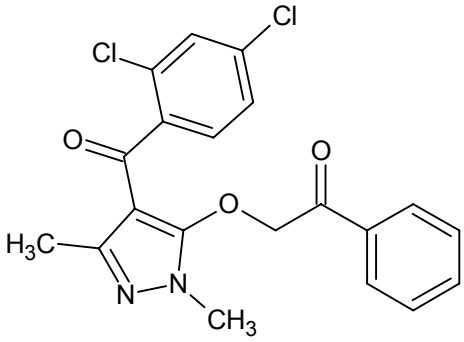
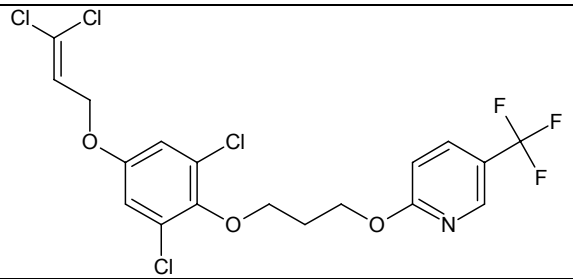
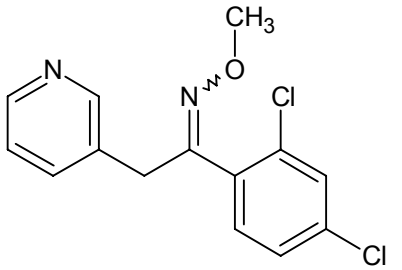
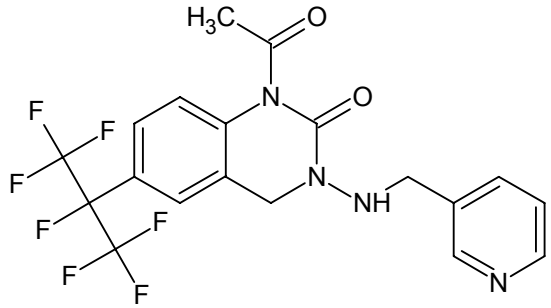
E pyrazoxyfen **F** pyrazoxyfène (m)	2-[4-(2,4-dichlorobenzoyl)-1,3-dimethylpyrazol-5-yloxy]acetophenone 2-[4-(2,4-dichlorobenzoyl)-1,3-diméthylpyrazol-5-yloxy]-1-phényléthanone 2-[[4-(2,4-dichlorobenzoyl)-1,3-dimethyl-1*H*-pyrazol-5-yl]oxy]-1-phenylethanone			**H**
		$C_{20}H_{16}Cl_2N_2O_3$	71561-11-0	
	InChI=1/C20H16Cl2N2O3/c1-12-18(19(26)15-9-8-14(21)10-16(15)22)20(24(2)23-12)27-11-17(25)13-6-4-3-5-7-13/h3-10H,11H2,1-2H3			
E pyridalyl **F** pyridalyle (m)	2,6-dichloro-4-(3,3-dichloroallyloxy)phenyl 3-[5-(trifluoromethyl)-2-pyridyloxy]propyl ether 2-(3-((2,6-dichloro-4-(3,3-dichloroprop-2-ényloxy)phénoxy))-propoxy)-5-trifluorométhylpyridin 2-[3-[2,6-dichloro-4-[(3,3-dichloro-2-propenyl)oxy]phenoxy]propoxy]-5-(trifluoromethyl)pyridine			**I**
		$C_{18}H_{14}Cl_4F_3NO_3$	179101-81-6	
	InChI=1/C18H14Cl4F3NO3/c19-13-8-12(27-7-4-15(21)22)9-14(20)17(13)29-6-1-5-28-16-3-2-11(10-26-16)18(23,24)25/h2-4,8-10H,1,5-7H2			
E pyrifenox **F** pyrifénox (m)	2′,4′-dichloro-2-(3-pyridyl)acetophenone (*EZ*)-*O*-methyloxime 1-(2,4-dichlorophényl)-2-(pyrid-3-yl)éthanone-*O*-méthyloxime 1-(2,4-dichlorophenyl)-2-(3-pyridinyl)ethanone *O*-methyloxime			**F**
		$C_{14}H_{12}Cl_2N_2O$	88283-41-4	
	InChI=1/C14H12Cl2N2O/c1-19-18-14(7-10-3-2-6-17-9-10)12-5-4-11(15)8-13(12)16/h2-6,8-9H,7H2,1H3/b18-14?			
E pyrifluquinazon **F** pyrifluquinazone (f)	1-acetyl-1,2,3,4-tetrahydro-3-[(3-pyridylmethyl)amino]-6-[1,2,2,2-tetrafluoro-1-(trifluoromethyl)ethyl]quinazolin-2-one 1-acétyl-1,2,3,4-tétrahydro-3-[(3-pyridylméthyl)amino]-6-[1,2,2,2-tétrafluoro-1-(trifluorométhyl)éthyl]quinazolin-2-one 1-acetyl-3,4-dihydro-3-[(3-pyridinylmethyl)amino]-6-[1,2,2,2-tetrafluoro-1-(trifluoromethyl)ethyl]-2(1*H*)-quinazolinone			**I**
		$C_{19}H_{15}F_7N_4O_2$	337458-27-2	
	InChI=1/C19H15F7N4O2/c1-11(31)30-15-5-4-14(17(20,18(21,22)23)19(24,25)26)7-13(15)10-29(16(30)32)28-9-12-3-2-6-27-8-12/h2-8,28H,9-10H2,1H3			

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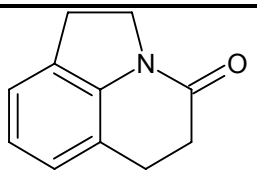
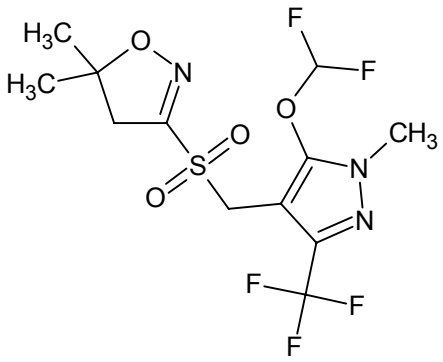
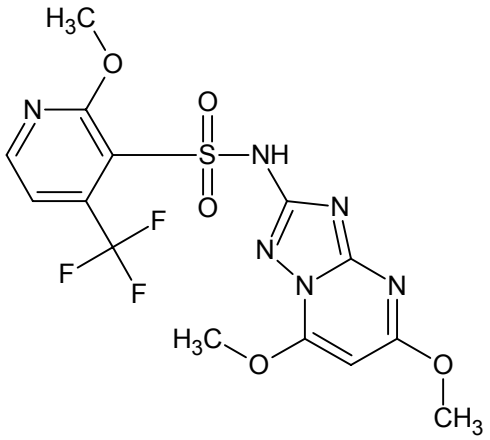
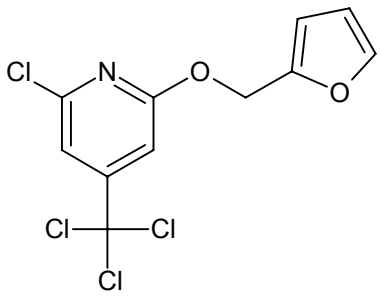
E pyroquilon **F** pyroquilone (f)	1,2,5,6-tetrahydropyrrolo[3,2,1-*ij*]quinolin-4-one 1,2,5,6-tétrahydropyrrolo[3,2,1-*ij*]quinolin-4-one 1,2,5,6-tetrahydro-4*H*-pyrrolo[3,2,1-*ij*]quinolin-4-one			**F**
		$C_{11}H_{11}NO$	57369-32-1	
	InChI=1/C11H11NO/c13-10-5-4-8-2-1-3-9-6-7-12(10)11(8)9/h1-3H,4-7H2			
E pyroxasulfone **F** pyroxasulfone (m)	3-[5-(difluoromethoxy)-1-methyl-3-(trifluoromethyl)pyrazol-4-ylmethylsulfonyl]-4,5-dihydro-5,5-dimethyl-1,2-oxazole 3-[(5-difluorométhoxy-1-méthyl-3-(trifluorométhyl)pyrazol-4-yl)méthyl-sulfonyl]-4,5-dihydro-5,5-diméthylisoxazole 3-[[[5-(difluoromethoxy)-1-methyl-3-(trifluoromethyl)-1*H*-pyrazol-4-yl]methyl]sulfonyl]-4,5-dihydro-5,5-dimethylisoxazole			**H**
		$C_{12}H_{14}F_5N_3O_4S$	447399-55-5	
	InChI=1/C12H14F5N3O4S/c1-11(2)4-7(19-24-11)25(21,22)5-6-8(12(15,16)17)18-20(3)9(6)23-10(13)14/h10H,4-5H2,1-3H3			
E pyroxsulam **F** pyroxsulame (m)	*N*-(5,7-dimethoxy[1,2,4]triazolo[1,5-*a*]pyrimidin-2-yl)-2-methoxy-4-(trifluoromethyl)pyridine-3-sulfonamide *N*-(5,7-diméthoxy[1,2,4]triazolo[1,5-*a*]pyrimidin-2-yl)-2-méthoxy-4-(trifluorométhyl)pyridine-3-sulfonamide *N*-(5,7-dimethoxy[1,2,4]triazolo[1,5-*a*]pyrimidin-2-yl)-2-methoxy-4-(trifluoromethyl)-3-pyridinesulfonamide			**H**
		$C_{14}H_{13}F_3N_6O_5S$	422556-08-9	
	InChI=1/C14H13F3N6O5S/c1-26-8-6-9(27-2)23-13(19-8)20-12(21-23)22-29(24,25)10-7(14(15,16)17)4-5-18-11(10)28-3/h4-6H,1-3H3,(H,21,22)/f/h22H			
E pyroxyfur **F** pyroxyfur (m)	6-chloro-4-trichloromethyl-2-pyridyl furfuryl ether 2-chloro-6-furfuryloxy-4-trichlorométhylpyridine 2-chloro-6-(2-furanylmethoxy)-4-(trichloromethyl)pyridine			**F**
		$C_{11}H_7Cl_4NO_2$	70166-48-2	
	InChI=1/C11H7Cl4NO2/c12-9-4-7(11(13,14)15)5-10(16-9)18-6-8-2-1-3-17-8/h1-5H,6H2			

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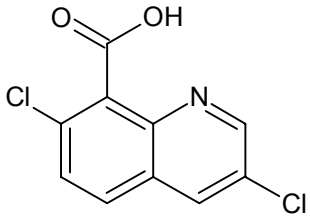
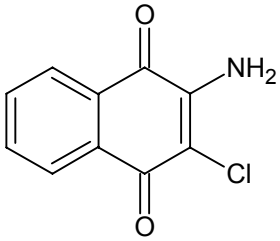
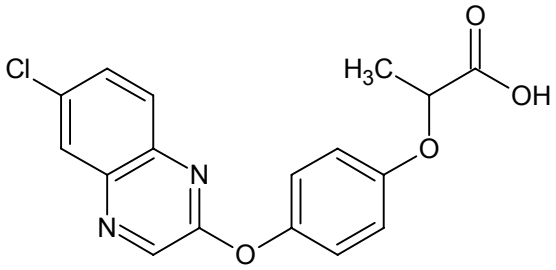
E quinclorac **F** quinclorac (m)	3,7-dichloroquinoline-8-carboxylic acid acide 3,7-dichloroquinoléine-8-carboxylique 3,7-dichloro-8-quinolinecarboxylic acid			**H**
		$C_{10}H_5Cl_2NO_2$	84087-01-4	
	InChI=1/C10H5Cl2NO2/c11-6-3-5-1-2-7(12)8(10(14)15)9(5)13-4-6/h1-4H,(H,14,15)/f/h14H			
E quinoclamine **F** quinoclamine (f)	2-amino-3-chloro-1,4-naphthoquinone 2-amino-3-chloro-1,4-naphtoquinone 2-amino-3-chloro-1,4-naphthalenedione			**AL** **H**
		$C_{10}H_6ClNO_2$	2797-51-5	
	InChI=1/C10H6ClNO2/c11-7-8(12)10(14)6-4-2-1-3-5(6)9(7)13/h1-4H,12H2			
E quizalofop **F** quizalofop (m)	(*RS*)-2-[4-(6-chloroquinoxalin-2-yloxy)phenoxy]propionic acid acide (*RS*)-2-[4-(6-chloroquinoxalin-2-yloxy)phénoxy]propionique 2-[4-[(6-chloro-2-quinoxalinyl)oxy]phenoxy]propanoic acid			**H**
		$C_{17}H_{13}ClN_2O_4$	76578-12-6	
	InChI=1/C17H13ClN2O4/c1-10(17(21)22)23-12-3-5-13(6-4-12)24-16-9-19-15-8-11(18)2-7-14(15)20-16/h2-10H,1H3,(H,21,22)/t10-/s3/f/h21H quizalofop-ethyl InChI=1/C19H17ClN2O4/c1-3-24-19(23)12(2)25-14-5-7-15(8-6-14)26-18-11-21-17-10-13(20)4-9-16(17)22-18/h4-12H,3H2,1-2H3/t12-/s3			
	NOTE It should be stated which ester or salt is present, for example quizalofop-ethyl [76578-14-8]. NOTE Il convient de préciser quel est l'ester ou le sel présent, par exemple quizalofop-éthyle [76578-14-8].			

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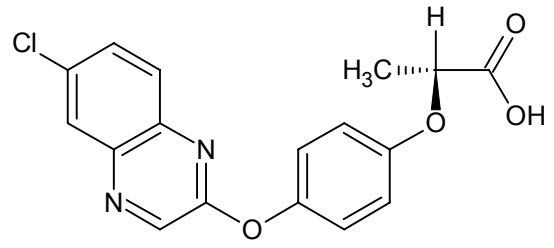
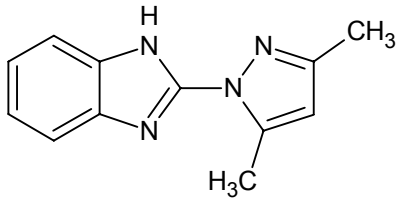
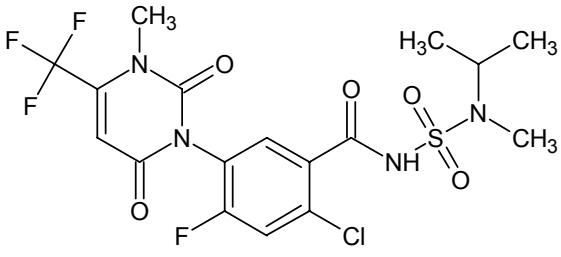
E quizalofop-P **F** quizalofop-P (m)	(*R*)-2-[4-(6-chloroquinoxalin-2-yloxy)phenoxy]propionic acid acide (*R*)-2-[4-(6-chloroquinoxalin-2-yloxy)phénoxy]propionique (2*R*)-2-[4-[(6-chloro-2-quinoxalinyl)oxy]phenoxy]propanoic acid	 $C_{17}H_{13}ClN_2O_4$	94051-08-8	**H**
	InChI=1/C17H13ClN2O4/c1-10(17(21)22)23-12-3-5-13(6-4-12)24-16-9-19-15-8-11(18)2-7-14(15)20-16/h2-10H,1H3,(H,21,22)/t10-/m1/s1/f/h21H quizalofop-P-ethyl InChI=1/C19H17ClN2O4/c1-3-24-19(23)12(2)25-14-5-7-15(8-6-14)26-18-11-21-17-10-13(20)4-9-16(17)22-18/h4-12H,3H2,1-2H3/t12-/m1/s1 quizalofop-P-tefuryl InChI=1/C22H21ClN2O5/c1-14(22(26)28-13-18-3-2-10-27-18)29-16-5-7-17(8-6-16)30-21-12-24-20-11-15(23)4-9-19(20)25-21/h4-9,11-12,14,18H,2-3,10,13H2,1H3/t14-,18?/m1/s1			
	NOTE It should be stated which ester or salt is present, for example quizalofop-P-ethyl [100646-51-3] or quizalofop-P-tefuryl [119738-06-6]. NOTE Il convient de préciser quel est l'ester ou le sel présent, par exemple quizalofop-P-éthyle [100646-51-3] ou quizalofop-P-tefuryl [119738-06-6].			
E rabenzazole **F** rabenzazole (m)	2-(3,5-dimethylpyrazol-1-yl)benzimidazole 2-(3,5-diméthyl-1*H*-pyrazol-1-yl)benzimidazole 2-(3,5-dimethyl-1*H*-pyrazol-1-yl)-1*H*-benzimidazole	 $C_{12}H_{12}N_4$	40341-04-6	**F**
	InChI=1/C12H12N4/c1-8-7-9(2)16(15-8)12-13-10-5-3-4-6-11(10)14-12/h3-7H,1-2H3,(H,13,14)/f/h13H			
E saflufenacil **F** saflufénacile (m)	*N'*-{2-chloro-4-fluoro-5-[1,2,3,6-tetrahydro-3-methyl-2,6-dioxo-4-(trifluoromethyl)pyrimidin-1-yl]benzoyl}-*N*-isopropyl-*N*-methylsulfamide *N'*-[2-chloro-4-fluoro-5-(3-méthyl-2,6-dioxo-4-(trifluorométhyl)-3,6-dihydro-1(2*H*)-pyrimidinyl)benzoyl]-*N*-isopropyl-*N*-méthylsulfamide 2-chloro-5-[3,6-dihydro-3-methyl-2,6-dioxo-4-(trifluoromethyl)-1(2*H*)-pyrimidinyl]-4-fluoro-*N*-[[methyl(1-methylethyl)amino]sulfonyl]benzamide	 $C_{17}H_{17}ClF_4N_4O_5S$	372137-35-4	**H**
	nChI=1/C17H17ClF4N4O5S/c1-8(2)25(4)32(30,31)23-15(28)9-5-12(11(19)6-10(9)18)26-14(27)7-13(17(20,21)22)24(3)16(26)29/h5-8H,1-4H3,(H,23,28)/f/h23H			

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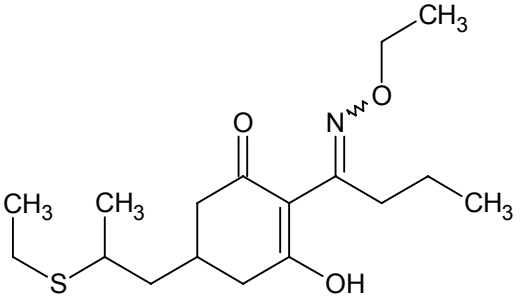
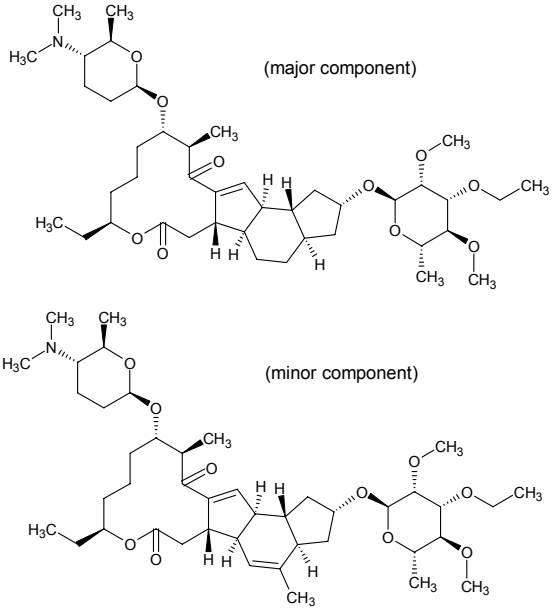
E sethoxydim **F** séthoxydime (f)	(5*RS*)-2-[(*EZ*)-1-(ethoxyimino)butyl]-5-[(2*RS*)-2-(ethylthio)propyl]-3-hydroxycyclohex-2-en-1-one (5*RS*)-2-[(*EZ*)-1-(éthoxyimino)butyl]-5-[(2*RS*)-2-(éthylthio)propyl]-3-hydroxycyclohex-2-èn-1-one 2-[1-(ethoxyimino)butyl]-5-[2-(ethylthio)propyl]-3-hydroxy-2-cyclohexen-1-one			**H**
		$C_{17}H_{29}NO_3S$	74051-80-2	
	InChI=1/C17H29NO3S/c1-5-8-14(18-21-6-2)17-15(19)10-13(11-16(17)20)9-12(4)22-7-3/h12-13,19H,5-11H2,1-4H3/b18-14?/t12?,13?			
E spinetoram **F** spinétorame (f)	mixture of 50 % to 90 % (1*S*,2*R*,5*R*,7*R*,9*R*,10*S*,14*R*,15*S*,19*S*)-7-(6-deoxy-3-*O*-ethyl-2,4-di-*O*-methyl-α-L-mannopyranosyloxy)-15-[(2*R*,5*S*,6*R*)-5-(dimethylamino)=tetrahydro-6-methylpyran-2-yloxy]-19-ethyl-14-methyl-20-oxa=tetracyclo[10.10.0.0^{2,10}.0^{5,9}]docos-11-ene-13,21-dione and 50 % to 10 % (1*S*,2*S*,5*R*,7*R*,9*S*,10*S*,14*R*,15*S*,19*S*)-7-(6-deoxy-3-*O*-ethyl-2,4-di-*O*-methyl-α-L-mannopyranosyloxy)-15-[(2*R*,5*S*,6*R*)-5-(dimethylamino)=tetrahydro-6-methylpyran-2-yloxy]-19-ethyl-4,14-dimethyl-20-oxa=tetracyclo[10.10.0.0^{2,10}.0^{5,9}]docosa-3,11-diene-13,21-dione mélange de (2*R*,3a*R*,5a*R*,5b*S*,9*S*,13*S*,14*R*,16a*S*,=16b*R*)-2-(6-désoxy-3-éthoxy-2,4-diméthoxy-α-L-mannopyranosyloxy)-13-[(2*R*,5*S*,6*R*)-5-(diméthylamino)=tétrahydro-6-méthylpyran-2-yloxy]-9-éthyl-2,3,3a,4,5,5a,5b,6,9,10,11,12,=13,14,16a,16b-hexadécahydro-14-méthyl-1*H*-*as*-indacéno=[3,2-*d*]oxacyclododécine-7,15-dione et de : (2*S*,3a*R*,5a*S*,5b*S*,9*S*,13*S*,14*R*,16a*S*,=16b*S*)-2-(6-désoxy-3-éthoxy-2,4-diméthoxy-α-L-mannopyranosyloxy)-13-[(2*R*,5*S*,6*R*)-5-(diméthylamino)=tétrahydro-6-méthylpyran-2-yloxy]-9-éthyl-2,3,3a,5a,5b,6,9,10,11,12,13,=14,16a,16b-tétradécahydro-4,14-diméthyl-1*H*-*as*-indacéno=[3,2-*d*]oxacyclododécine-7,15-dione dans les proportions de 50–90% et 50–10%			**I**

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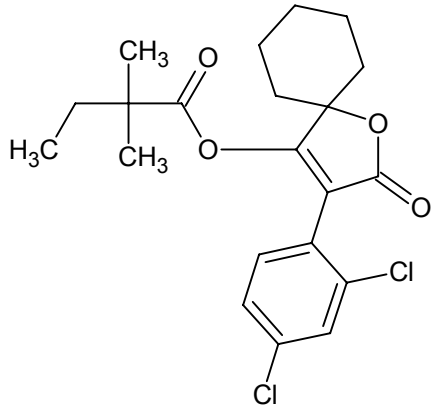
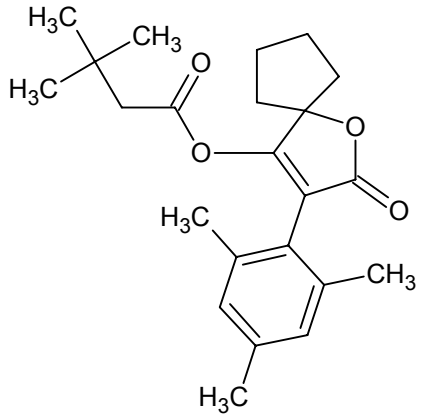
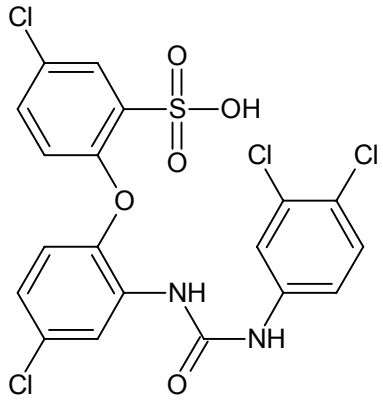
	IUPAC International Chemical Identifier (InChI™)			
	(2*R*,3a*R*,5a*R*,5b*S*,9*S*,13*S*,14*R*,16a*S*,= 16b*R*)-2-[(6-deoxy-3-*O*-ethyl- 2,4-di-*O*-methyl-α-L-mannopyranosyl)= oxy]-13-[[(2*R*,5*S*,6*R*)-5-(dimethyl= amino)tetrahydro-6-methyl- 2*H*-pyran-2-yl]oxy]-9-ethyl- 2,3,3a,4,5,5a,5b,6,9,10,11,12,13,14,= 16a,16b-hexadecahydro-14-methyl- 1*H*-*as*-indaceno[3,2-*d*]oxacyclo= dodecin-7,15-dione mixture with (2*S*,3a*R*,5a*S*,5b*S*,9*S*,13*S*,14*R*,16a*S*,= 16b*S*)-2-[(6-deoxy-3-*O*-ethyl- 2,4-di-*O*-methyl-α-L-mannopyranosyl)= oxy]-13-[[(2*R*,5*S*,6*R*)-5-(dimethyl= amino)tetrahydro-6-methyl- 2*H*-pyran-2-yl]oxy]-9-ethyl- 2,3,3a,5a,5b,6,9,10,11,12,13,14,= 16a,16b-tetradecahydro-4,14-dimethyl- 1*H*-*as*-indaceno[3,2-*d*]oxacyclo= dodecin-7,15-dione	$C_{42}H_{69}NO_{10}$ + $C_{43}H_{69}NO_{10}$	187166-40-1 + 187166-15-0	
	major component (4,5-dihydro) InChI=1/C42H69NO10/c1-10-27-13-12-14-35(53-37-18-17-34(43(6)7)24(4)49-37)23(3)38(45)33-21-31-29(32(33)22-36(44)51-27)16-15-26-19-28(20-30(26)31)52-42-41(47-9)40(48-11-2)39(46-8)25(5)50-42/h21,23-32,34-35,37,39-42H,10-20,22H2,1-9H3/t23-,24-,25+,26-,27+,28-,29-,30-,31-,32+,34+,35+,37+,39+,40-,41-,42+/m1/s1 minor component (4-methyl) InChI=1/C43H69NO10/c1-11-27-14-13-15-36(54-38-17-16-35(44(7)8)25(5)50-38)24(4)39(46)34-21-32-30(33(34)22-37(45)52-27)18-23(3)29-19-28(20-31(29)32)53-43-42(48-10)41(49-12-2)40(47-9)26(6)51-43/h18,21,24-33,35-36,38,40-43H,11-17,19-20,22H2,1-10H3/t24-,25-,26+,27+,28-,29+,30-,31-,32-,33+,35+,36+,38+,40+,41-,42-,43+/m1/s1			
E spirodiclofen **F** spirodiclofène (m)	3-(2,4-dichlorophenyl)-2-oxo-1-oxaspiro[4.5]dec-3-en-4-yl 2,2-dimethylbutyrate 2,2-diméthylbutyrate de 3-(2,4-dichlorophényl)-2-oxo-1-oxaspiro[4.5]déc-3-èn-4-yle 3-(2,4-dichlorophenyl)-2-oxo-1-oxaspiro[4.5]dec-3-en-4-yl 2,2-dimethylbutanoate			**A**
		$C_{21}H_{24}Cl_2O_4$	148477-71-8	
	InChI=1/C21H24Cl2O4/c1-4-20(2,3)19(25)26-17-16(14-9-8-13(22)12-15(14)23)18(24)27-21(17)10-6-5-7-11-21/h8-9,12H,4-7,10-11H2,1-3H3			

E: Common name F: Nom commun	Chemical name Nom chimique E: IUPAC F: UICPA C: CAS	Structure Structure Molecular formula Formule brute	CAS Registry Number® Numéro d'enregistrement CAS®	Use Applic-ation
	IUPAC International Chemical Identifier (InChI™)			
E spiromesifen **F** spiromésifène (m)	3-mesityl-2-oxo-1-oxaspiro[4.4]non-3-en-4-yl 3,3-dimethylbutyrate 3,3-diméthylbutyrate de 3-mésityl-2-oxo-1-oxaspiro[4.4]non-3-èn-4-yle 2-oxo-3-(2,4,6-trimethylphenyl)-1-oxaspiro[4.4]non-3-en-4-yl 3,3-dimethylbutanoate			I
		$C_{23}H_{30}O_4$	283594-90-1	
	InChI=1/C23H30O4/c1-14-11-15(2)18(16(3)12-14)19-20(26-17(24)13-22(4,5)6)23(27-21(19)25)9-7-8-10-23/h11-12H,7-10,13H2,1-6H3			
E sulcofuron **F** sulcofuron (m)	5-chloro-2-{4-chloro-2-[3-(3,4-dichlorophenyl)ureido]phenoxy}= benzenesulfonic acid acide 5-chloro-2-[4-chloro-2-[3-(3,4-dichlorophényl)uréido]phénoxy]= benzène-sulfonique 5-chloro-2-[4-chloro-2-[[[(3,4-dichlorophenyl)amino]carbonyl]= amino]phenoxy]benzenesulfonic acid			I
		$C_{19}H_{12}Cl_4N_2O_5S$	24019-05-4	
	InChI=1/C19H12Cl4N2O5S/c20-10-1-5-16(30-17-6-2-11(21)8-18(17)31(27,28)29)15(7-10)25-19(26)24-12-3-4-13(22)14(23)9-12/h1-9H,(H2,24,25,26)(H,27,28,29)/f/h24-25,27H sulcofuron-sodium InChI=1/C19H12Cl4N2O5S.Na/c20-10-1-5-16(30-17-6-2-11(21)8-18(17)31(27,28)29)15(7-10)25-19(26)24-12-3-4-13(22)14(23)9-12;/h1-9H,(H2,24,25,26)(H,27,28,29);/q;+1/p-1/fC19H11Cl4N2O5S.Na/h24-25H;/q-1;m			
	NOTE It should be stated which ester or salt is present, for example sulcofuron-sodium [3567-25-7]. NOTE Il convient de préciser quel est l'ester ou le sel présent, par exemple sulcofuron-sodium [3567-25-7].			

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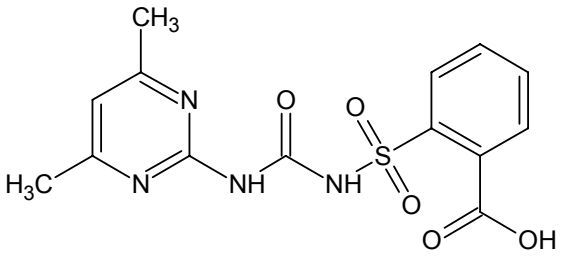
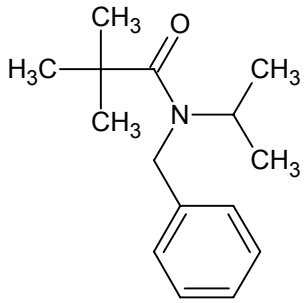
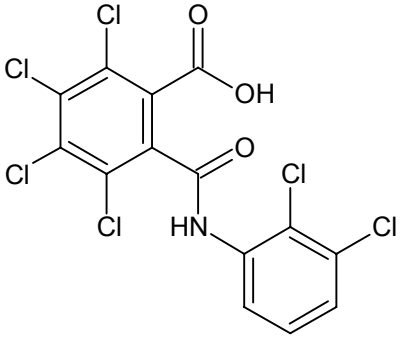
E sulfometuron **F** sulfométuron (m)	2-(4,6-dimethylpyrimidin-2-ylcarbamoylsulfamoyl)benzoic acid acide *N*-(4,6-diméthylpyrimidin-2-yl-carbamyl)-*o*-sulfaminobenzoïque 2-[[[[(4,6-dimethyl-2-pyrimidinyl)=amino]carbonyl]amino]sulfonyl]=benzoic acid			**H**
		$C_{14}H_{14}N_4O_5S$	74223-56-6	
	InChI=1/C14H14N4O5S/c1-8-7-9(2)16-13(15-8)17-14(21)18-24(22,23)11-6-4-3-5-10(11)12(19)20/h3-7H,1-2H3,(H,19,20)(H2,15,16,17,18,21)/f/h17-19H sulfometuron-methyl InChI=1/C15H16N4O5S/c1-9-8-10(2)17-14(16-9)18-15(21)19-25(22,23)12-7-5-4-6-11(12)13(20)24-3/h4-8H,1-3H3,(H2,16,17,18,19,21)/f/h18-19H			
	NOTE It should be stated which ester or salt is present, for example sulfometuron-methyl [74222-97-2]. NOTE Il convient de préciser quel est l'ester ou le sel présent, par exemple sulfométuron-méthyle [74222-97-2].			
E tebutam **F** tébutame (m)	*N*-benzyl-*N*-isopropyl-2,2-dimethylpropionamide *N*-benzyl-*N*-isopropylpivalamide 2,2-dimethyl-*N*-(1-methylethyl)-*N*-(phenylmethyl)propanamide			**H**
		$C_{15}H_{23}NO$	35256-85-0	
	InChI=1/C15H23NO/c1-12(2)16(14(17)15(3,4)5)11-13-9-7-6-8-10-13/h6-10,12H,11H2,1-5H3			
	NOTE 1 The name “tebutam” is not acceptable for use in the USA, where “butam” has been adopted as the common name. NOTE 1 Le nom «tebutam» n'est pas acceptable pour l'emploi aux États-Unis, où «butam» a été adopté comme nom commun. NOTE 2 The name “tebutam” is not acceptable for use in Japan because of the risk of confusion with the trade name “Tepdum”. NOTE 2 Le nom «tebutam» n'est pas acceptable pour l'emploi au Japon, car il entre en conflit avec le nom commercial «Tepdum».			
E tecloftalam **F** técloftalame (m)	2′,3,3′,4,5,6-hexachlorophthalanilic acid acide 2′,3,3′,4,5,6–hexachlorophtalanilique 2,3,4,5-tetrachloro-6-[[(2,3-dichlorophenyl)amino]carbonyl]=benzoic acid			**B** **F**
		$C_{14}H_5Cl_6NO_3$	76280-91-6	
	InChI=1/C14H5Cl6NO3/c15-4-2-1-3-5(8(4)16)21-13(22)6-7(14(23)24)10(18)12(20)11(19)9(6)17/h1-3H,(H,21,22)(H,23,24)/f/h21,23H			

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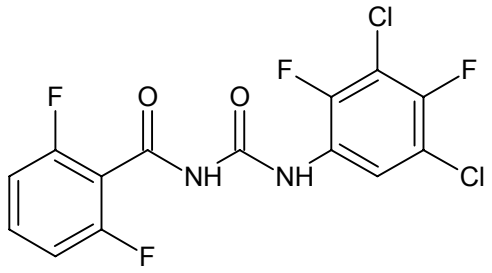
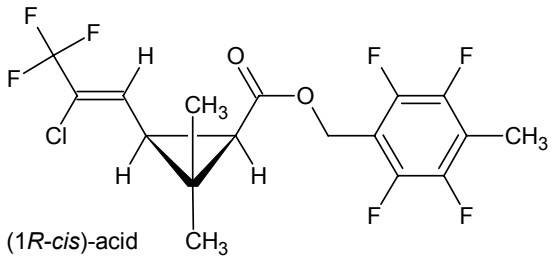
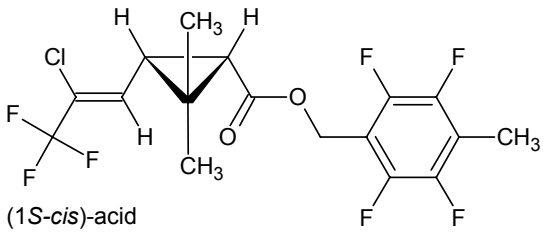
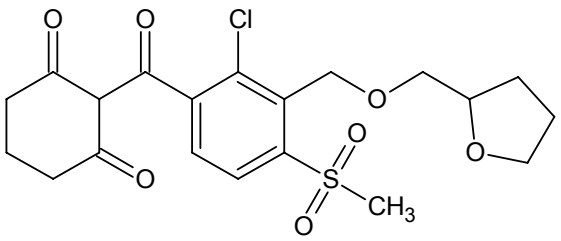
E teflubenzuron **F** téflubenzuron (m)	1-(3,5-dichloro-2,4-difluorophenyl)-3-(2,6-difluorobenzoyl)urea *N*-(3,5-dichloro-2,4-difluorophényl)-*N'*-(2,6-difluorobenzoyl)urée ou 1-(3,5-dichloro-2,4-difluorophényl)-3-(2,6-difluorobenzoyl)urée *N*-[[(3,5-dichloro-2,4-difluorophenyl)amino]carbonyl]-2,6-difluorobenzamide	 $C_{14}H_6Cl_2F_4N_2O_2$	83121-18-0	**IGR**
	InChI=1/C14H6Cl2F4N2O2/c15-5-4-8(12(20)10(16)11(5)19)21-14(24)22-13(23)9-6(17)2-1-3-7(9)18/h1-4H,(H2,21,22,23,24)/f/h21-22H			
E tefluthrin **F** téfluthrine (f)	2,3,5,6-tetrafluoro-4-methylbenzyl (1*RS*,3*RS*)-3-[(*Z*)-2-chloro-3,3,3-trifluoroprop-1-enyl]-2,2-dimethylcyclopropanecarboxylate Alternative Rothamsted-style stereodescriptors: 2,3,5,6-tetrafluoro-4-methylbenzyl (1*RS*)-*cis*-3-[(*Z*)-2-chloro-3,3,3-trifluoroprop-1-enyl]-2,2-dimethylcyclopropanecarboxylate (1*RS*,3*RS*)-((*Z*)-2-chloro-3,3,3-trifluoroprop-1-ènyl)-2,2-diméthylcyclopropanecarboxylate de 2,3,5,6-tétrafluoro-4-méthylbenzyle (2,3,5,6-tetrafluoro-4-methylphenyl)=methyl (1*R*,3*R*)-*rel*-3-[(1*Z*)-2-chloro-3,3,3-trifluoro-1-propenyl]-2,2-dimethylcyclopropanecarboxylate	 (1*R-cis*)-acid  (1*S-cis*)-acid $C_{17}H_{14}ClF_7O_2$	79538-32-2	**I**
	InChI=1/C17H14ClF7O2/c1-6-11(19)13(21)7(14(22)12(6)20)5-27-15(26)10-8(16(10,2)3)4-9(18)17(23,24)25/h4,8,10H,5H2,1-3H3/b9-4-/t8-,10-/s3			
E tefuryltrione **F** téfuryltrione (f)	2-{2-chloro-4-mesyl-3-[(*RS*)-tetrahydrofuran-2-ylmethoxymethyl]=benzoyl}cyclohexane-1,3-dione 2-[2-chloro-4-mésyl-3-(tétrahydrofuran-2-yl(méthoxy)méthyl)-benzoyl]cyclohexane-1,3-dione 2-[2-chloro-4-(methylsulfonyl)-3-[[(tetrahydro-2-furanyl)methoxy]=methyl]benzoyl]-1,3-cyclohexanedione	 $C_{20}H_{23}ClO_7S$	473278-76-1	**H**
	InChI=1/C20H23ClO7S/c1-29(25,26)17-8-7-13(20(24)18-15(22)5-2-6-16(18)23)19(21)14(17)11-27-10-12-4-3-9-28-12/h7-8,12,18H,2-6,9-11H2,1H3/t12-/s3			

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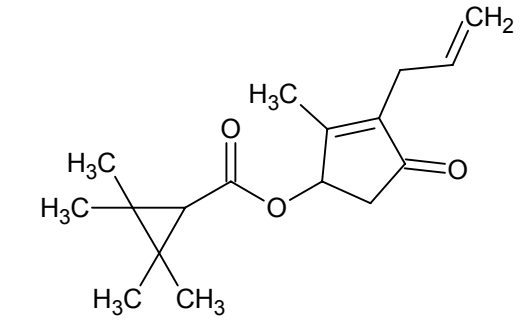
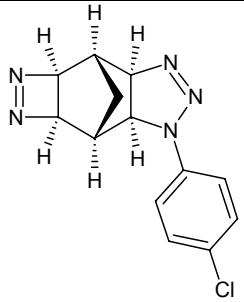
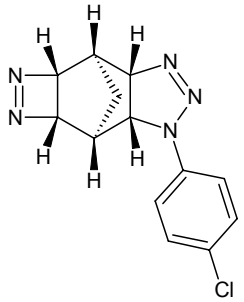
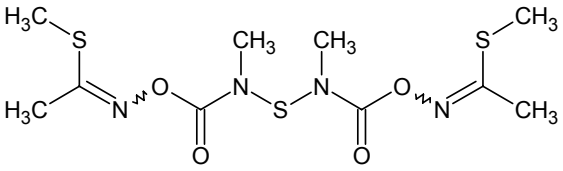
E terallethrin **F** tétralléthrine (f)	(*RS*)-3-allyl-2-methyl-4-oxocyclopent-2-enyl 2,2,3,3-tetramethylcyclopropanecarboxylate 2,2,3,3-tétraméthylcyclopropanecarboxylate de (*RS*)-3-allyl-2-méthyl-4-oxocyclopent-2-ènyle 2-methyl-4-oxo-3-(2-propenyl)-2-cyclopenten-1-yl 2,2,3,3-tetramethylcyclopropanecarboxylate			**I**
		$C_{17}H_{24}O_3$	15589-31-8	
	InChI=1/C17H24O3/c1-7-8-11-10(2)13(9-12(11)18)20-15(19)14-16(3,4)17(14,5)6/h7,13-14H,1,8-9H2,2-6H3/t13-/s3			
E tetcyclacis **F** tetcyclacis (m)	*rel*-(1*R*,2*R*,6*S*,7*R*,8*R*,11*S*)-5-(4-chlorophenyl)-3,4,5,9,10-pentaazatetracyclo[5.4.1.0^{2,6}.0^{8,11}]=dodeca-3,9-diene *rel*-(3a*R*,4*R*,4a*S*,6a*R*,7*R*,7a*S*)-1-(4-chlorophényl)-(3a,4,4a,6a,7,7a)-hexahydro-4,7-méthano-1*H*-[1,2-diazéto][3,4-*f*]benzotriazole (3a*R*,4*R*,4a*S*,6a*R*,7*R*,7a*S*)-*rel*-1-(4-chlorophenyl)-3a,4,4a,6a,7,7a-hexahydro-4,7-methano-1*H*-[1,2]diazeto[3,4-*f*]benzotriazole	 (1*R*,2*R*,6*S*,7*R*,8*R*,11*S*)-isomer  (1*S*,2*S*,6*R*,7*S*,8*S*,11*R*)-isomer		**P**
		$C_{13}H_{12}ClN_5$	77788-21-7	
	InChI=1/C13H12ClN5/c14-6-1-3-7(4-2-6)19-13-9-5-8(12(13)17-18-19)10-11(9)16-15-10/h1-4,8-13H,5H2/t8-,9+,10+,11-,12-,13+/s3			
E thiodicarb **F** thiodicarb (m)	(3*EZ*,12*EZ*)-3,7,9,13-tetramethyl-5,11-dioxa-2,8,14-trithia-4,7,9,12-tetraazapentadeca-3,12-diene-6,10-dione (3*EZ*,12*EZ*)-3,7,9,13-tétraméthyl-5,11-dioxa-2,8,14-trithia-4,7,9,12-tétra-azapentadéca-3,12-diène-6,10-dione dimethyl *N,N'*-[thiobis[(methylimino)=carbonyloxy]]bis[ethanimidothioate]			**I** **M**
		$C_{10}H_{18}N_4O_4S_3$	59669-26-0	
	InChI=1/C10H18N4O4S3/c1-7(19-5)11-17-9(15)13(3)21-14(4)10(16)18-12-8(2)20-6/h1-6H3/b11-7?,12-8?			

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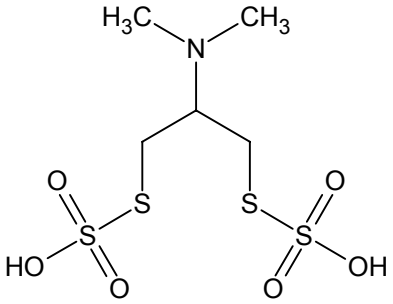
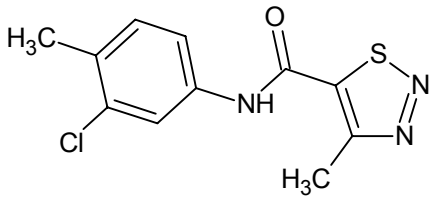
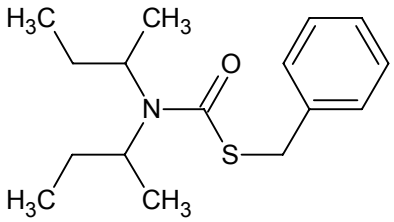
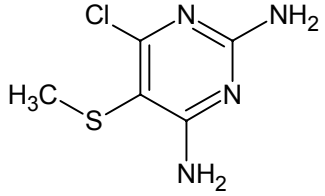
E thiosultap **F** thiosultap (m)	dihydrogen *S,S′*-[2-(dimethylamino)trimethylene]= di(thiosulfate) di(hydrogénothiosulfate) de 2-diméthylaminopropane-1,3-diyle thiosulfuric acid ($H_2S_2O_3$) *S,S′*-[2-(dimethylamino)-1,3-propanediyl] ester			**I**
		$C_5H_{13}NO_6S_4$	98968-92-4	
	InChI=1/C5H13NO6S4/c1-6(2)5(3-13-15(7,8)9)4-14-16(10,11)12/h5H,3-4H2,1-2H3,(H,7,8,9)(H,10,11,12)/f/h7,10H thiosultap-monosodium InChI=1/C5H13NO6S4.Na/c1-6(2)5(3-13-15(7,8)9)4-14-16(10,11)12;/h5H,3-4H2,1-2H3,(H,7,8,9)(H,10,11,12);/q;+1/p-1/fC5H12NO6S4.Na/h7H;/q-1;m thiosultap-disodium InChI=1/C5H13NO6S4.2Na/c1-6(2)5(3-13-15(7,8)9)4-14-16(10,11)12;;/h5H,3-4H2,1-2H3,(H,7,8,9)(H,10,11,12);;/q;2*+1/p-2/fC5H11NO6S4.2Na/q-2;2m			
	NOTE It should be stated which ester or salt is present, for example thiosultap-monosodium [29547-00-0] or thiosultap-disodium [52207-48-4]. NOTE Il convient de préciser quel est l'ester ou le sel présent, par exemple thiosultap-monosodium [29547-00-0] ou thiosultap-disodique [52207-48-4].			
E tiadinil **F** tiadinil (m)	3′-chloro-4,4′-dimethyl-1,2,3-thiadiazole-5-carboxanilide *N*-(3-chloro-4-méthylphényl)-4-méthyl-1,2,3-thiadiazole-5-carboxamide *N*-(3-chloro-4-methylphenyl)-4-methyl-1,2,3-thiadiazole-5-carboxamide			**F**
		$C_{11}H_{10}ClN_3OS$	223580-51-6	
	InChI=1/C11H10ClN3OS/c1-6-3-4-8(5-9(6)12)13-11(16)10-7(2)14-15-17-10/h3-5H,1-2H3,(H,13,16)/f/h13H			
E tiocarbazil **F** tiocarbazil (m)	*S*-benzyl di-*sec*-butyl(thiocarbamate) di-*sec*-butylthiocarbamate de *S*-benzyle *S*-(phenylmethyl) bis(1-methylpropyl)carbamothioate			**H**
		$C_{16}H_{25}NOS$	36756-79-3	
	InChI=1/C16H25NOS/c1-5-13(3)17(14(4)6-2)16(18)19-12-15-10-8-7-9-11-15/h7-11,13-14H,5-6,12H2,1-4H3			
E tioclorim **F** tioclorime (f)	6-chloro-5-methylthiopyrimidine-2,4-diamine 6-chloro-5-méthylthiopyrimidin-2,4-diyldiamine 6-chloro-5-(methylthio)-2,4-pyrimidinediamine			**H**
		$C_5H_7ClN_4S$	68925-41-7	
	InChI=1/C5H7ClN4S/c1-11-2-3(6)9-5(8)10-4(2)7/h1H3,(H4,7,8,9,10)/f/h7-8H2			

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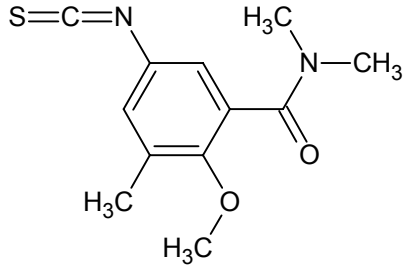
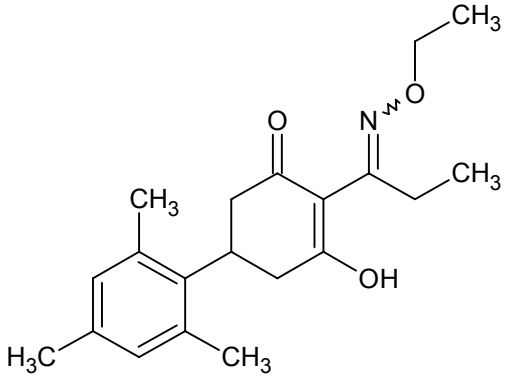
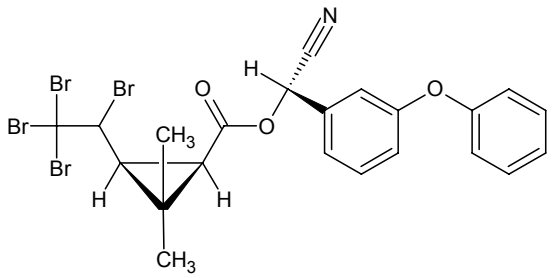
	IUPAC International Chemical Identifier (InChI™)			
E tioxymid **F** tioxymide (m)	5-isothiocyanato-*N,N*,3-trimethyl-*o*-anisamide 5-isothiocyanato-2-méthoxy-*N,N*,3-triméthylbenzamide 5-isothiocyanato-2-methoxy-*N,N*,3-trimethylbenzamide			**F**
		$C_{12}H_{14}N_2O_2S$	70751-94-9	
	InChI=1/C12H14N2O2S/c1-8-5-9(13-7-17)6-10(11(8)16-4)12(15)14(2)3/h5-6H,1-4H3			
E tralkoxydim **F** tralkoxydime (f)	(*RS*)-2-[(*EZ*)-1-(ethoxyimino)propyl]-3-hydroxy-5-mesitylcyclohex-2-en-1-one (*RS*)-2-[(*EZ*)-1-(éthoxyimino)propyl]-3-hydroxy-5-mésitylcyclohex-2-èn-1-one 2-[1-(ethoxyimino)propyl]-3-hydroxy-5-(2,4,6-trimethylphenyl)-2-cyclohexen-1-one			**H**
		$C_{20}H_{27}NO_3$	87820-88-0	
	InChI=1/C20H27NO3/c1-6-16(21-24-7-2)20-17(22)10-15(11-18(20)23)19-13(4)8-12(3)9-14(19)5/h8-9,15,22H,6-7,10-11H2,1-5H3/b21-16?/t15-/s3			
E tralomethrin **F** tralométhrine (f)	(*S*)-α-cyano-3-phenoxybenzyl (1*R*,3*S*)-2,2-dimethyl-3-[(*RS*)-1,2,2,2-tetrabromoethyl]cyclopropane= carboxylate Alternative Rothamsted-style stereodescriptors: (*S*)-α-cyano-3-phenoxybenzyl (1*R*)-*cis*-2,2-dimethyl-3-[(*RS*)-1,2,2,2-tetrabromoethyl]cyclopropane= carboxylate (1*R*,3*S*)-2,2-diméthyl-3-[(*RS*)-1,2,2,2-tétrabromoéthyl]cyclopropane= carboxylate de (*S*)-cyano-(3-phénoxyphényl)méthyle (*S*)-cyano(3-phenoxyphenyl)methyl (1*R*,3*S*)-2,2-dimethyl-3-(1,2,2,2-tetrabromoethyl)cyclopropane= carboxylate			**I**
		$C_{22}H_{19}Br_4NO_3$	66841-25-6	
	InChI=1/C22H19Br4NO3/c1-21(2)17(19(23)22(24,25)26)18(21)20(28)30-16(12-27)13-7-6-10-15(11-13)29-14-8-4-3-5-9-14/h3-11,16-19H,1-2H3/t16-,17-,18+,19?/m1/s1			
	NOTE An unequal mixture of the (*S*), (1*R*,3*S*), (*R*) and the (*S*), (1*R*,3*S*), (*S*) diastereoisomers. NOTE Un mélange inégal des diastéréoisomères (*S*), (1*R*,3*S*), (*R*) et (*S*), (1*R*,3*S*), (*S*).			

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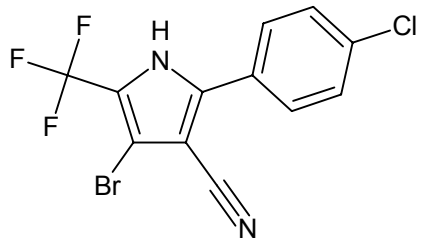
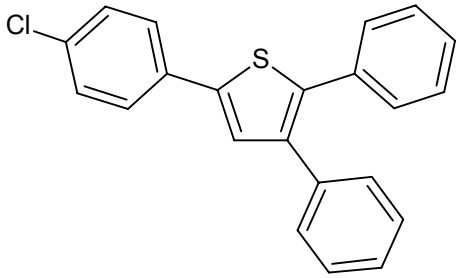
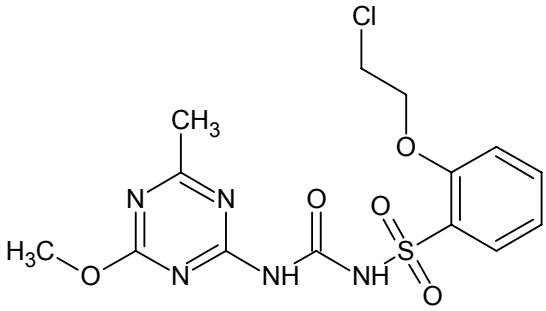
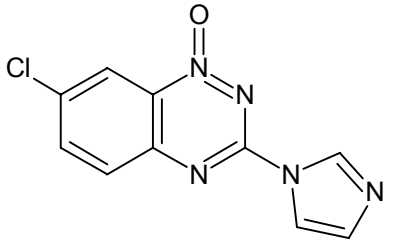
E tralopyril **F** tralopyril (m)	4-bromo-2-(4-chlorophenyl)-5-(trifluoromethyl)-1*H*-pyrrole-3-carbonitrile 4-bromo-2-(4-chlorophényl)-5-(trifluorométhyl)-1*H*-pyrrole-3-carbonitrile 4-bromo-2-(4-chlorophenyl)-5-(trifluoromethyl)-1*H*-pyrrole-3-carbonitrile	 $C_{12}H_5BrClF_3N_2$	122454-29-9	**M**
	InChI=1/C12H5BrClF3N2/c13-9-8(5-18)10(19-11(9)12(15,16)17)6-1-3-7(14)4-2-6/h1-4,19H			
E triarathene **F** triarathène (m)	5-(4-chlorophenyl)-2,3-diphenylthiophene 5-(4-chlorophényl)-2,3-diphénylthiophène 5-(4-chlorophenyl)-2,3-diphenylthiophene	 $C_{22}H_{15}ClS$	65691-00-1	**A** **I**
	InChI=1/C22H15ClS/c23-19-13-11-17(12-14-19)21-15-20(16-7-3-1-4-8-16)22(24-21)18-9-5-2-6-10-18/h1-15H			
E triasulfuron **F** triasulfuron (m)	1-[2-(2-chloroethoxy)phenylsulfonyl]-3-(4-methoxy-6-methyl-1,3,5-triazin-2-yl)urea *N*-[2-(2-chloroéthoxy)phénylsulfonyl]-*N'*-(4-méthoxy-6-méthyl-1,3,5-triazin-2-yl)urée ou 1-[2-(2-chloroéthoxy)phénylsulfonyl]-3-(4-méthoxy-6-méthyl-1,3,5-triazin-2-yl)urée 2-(2-chloroethoxy)-*N*-[[(4-methoxy-6-methyl-1,3,5-triazin-2-yl)amino]=carbonyl]benzenesulfonamide	 $C_{14}H_{16}ClN_5O_5S$	82097-50-5	**H**
	InChI=1/C14H16ClN5O5S/c1-9-16-12(19-14(17-9)24-2)18-13(21)20-26(22,23)11-6-4-3-5-10(11)25-8-7-15/h3-6H,7-8H2,1-2H3,(H2,16,17,18,19,20,21)/f/h18,20H			
E triazoxide **F** triazoxide (m)	7-chloro-3-imidazol-1-yl-1,2,4-benzotriazine 1-oxide 1-oxyde de 7-chloro-3-imidazol-1-yl-1,2,4-benzotriazine 7-chloro-3-(1*H*-imidazol-1-yl)-1,2,4-benzotriazine 1-oxide	 $C_{10}H_6ClN_5O$	72459-58-6	**F**
	InChI=1/C10H6ClN5O/c11-7-1-2-8-9(5-7)16(17)14-10(13-8)15-4-3-12-6-15/h1-6H			

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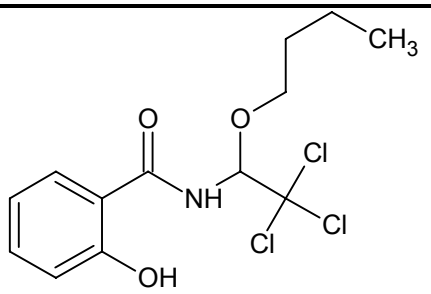
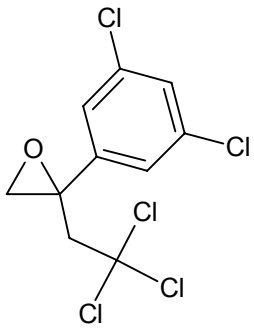
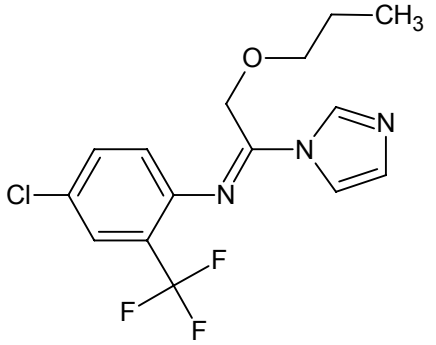
E trichlamide **F** trichlamide (m)	(*RS*)-*N*-(1-butoxy-2,2,2-trichloroethyl)salicylamide (*RS*)-*N*-(1-butoxy-2,2,2-trichloroéthyl)salicylamide *N*-(1-butoxy-2,2,2-trichloroethyl)-2-hydroxybenzamide			**F**
		$C_{13}H_{16}Cl_3NO_3$	70193-21-4	
	InChI=1/C13H16Cl3NO3/c1-2-3-8-20-12(13(14,15)16)17-11(19)9-6-4-5-7-10(9)18/h4-7,12,18H,2-3,8H2,1H3,(H,17,19)/t12-/s3/f/h17H			
E tridiphane **F** tridifane (m)	(*RS*)-2-(3,5-dichlorophenyl)-2-(2,2,2-trichloroethyl)oxirane (*RS*)-2-(3,5-dichlorophényl)-2-(2,2,2-trichloroéthyl)oxirane 2-(3,5-dichlorophenyl)-2-(2,2,2-trichloroethyl)oxirane			**H**
		$C_{10}H_7Cl_5O$	58138-08-2	
	InChI=1/C10H7Cl5O/c11-7-1-6(2-8(12)3-7)9(5-16-9)4-10(13,14)15/h1-3H,4-5H2/t9-/s3			
E triflumizole **F** triflumizole (f)	(*E*)-4-chloro-α,α,α-trifluoro-*N*-(1-imidazol-1-yl-2-propoxyethylidene)-*o*-toluidine 1-[(1*E*)-1-[[4-chloro-2-(trifluorométhyl)phényl]imino]-2-propoxyéthyl]-1*H*-imidazole 1-[(1*E*)-1-[[4-chloro-2-(trifluoromethyl)phenyl]imino]-2-propoxyethyl]-1*H*-imidazole			**F**
		$C_{15}H_{15}ClF_3N_3O$	99387-89-0	
	InChI=1/C15H15ClF3N3O/c1-2-7-23-9-14(22-6-5-20-10-22)21-13-4-3-11(16)8-12(13)15(17,18)19/h3-6,8,10H,2,7,9H2,1H3/b21-14+			

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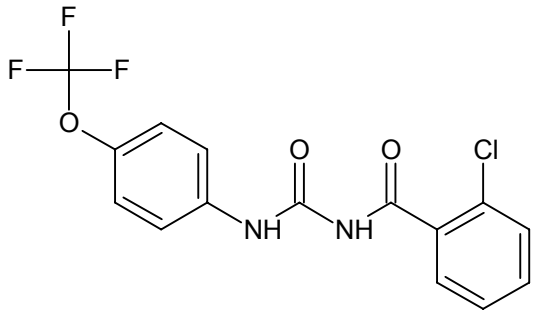
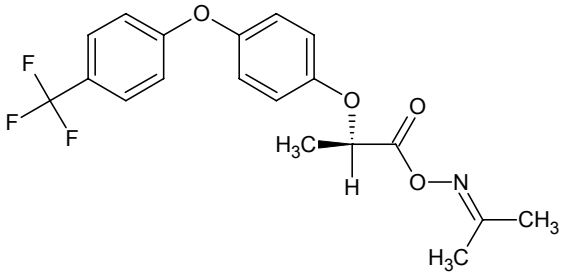
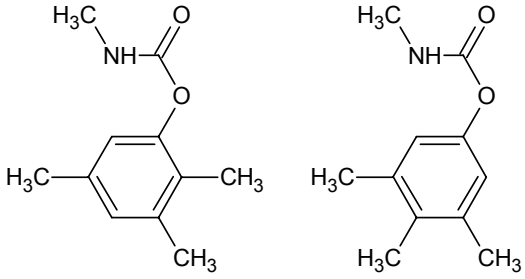
E triflumuron **F** triflumuron (m)	1-(2-chlorobenzoyl)-3-(4-trifluoromethoxyphenyl)urea *N*-(2-chlorobenzoyl)-*N'*-(4-trifluorométhoxyphényl)urée ou 1-(2-chlorobenzoyl)-3-(4-trifluorométhoxyphényl)urée 2-chloro-*N*-[[[4-(trifluoromethoxy)=phenyl]amino]carbonyl]benzamide	 $C_{15}H_{10}ClF_3N_2O_3$	64628-44-0	**IGR**
	InChI=1/C15H10ClF3N2O3/c16-12-4-2-1-3-11(12)13(22)21-14(23)20-9-5-7-10(8-6-9)24-15(17,18)19/h1-8H,(H2,20,21,22,23)/f/h20-21H			
E trifopsime **F** trifopsime (f)	acetone (*R*)-*O*-{2-[4-(α,α,α-trifluoro-*p*-tolyloxy)phenoxy]propionyl}oxime acétone-(*R*)-*O*-{2-[4-(4-trifluorométhylphénoxy)phénoxy]=propionyl}-oxime 2-propanone *O*-[(2*R*)-1-oxo-2-[4-[4-(trifluoromethyl)phenoxy]phenoxy]=propyl]oxime	 $C_{19}H_{18}F_3NO_4$	72131-76-1	**H**
	InChI=1/C19H18F3NO4/c1-12(2)23-27-18(24)13(3)25-15-8-10-17(11-9-15)26-16-6-4-14(5-7-16)19(20,21)22/h4-11,13H,1-3H3/t13-/m1/s1			
E trimethacarb **F** trimétacarb (m)	a reaction product comprising from 3,5 to 5 parts by mass of 3,4,5-trimethylphenyl methylcarbamate to 1 part by mass of 2,3,5-trimethylphenyl methylcarbamate méthylcarbamate de 2,3,5(ou 3,4,5)-triméthylphényle mélange comprenant de 3,5 à 5 parties d'isomère 3,4,5 pour une partie d'isomère 2,3,5 2,3,5(*or* 3,4,5)-trimethylphenyl methylcarbamate	 (2,3,5-isomer) (3,4,5-isomer) $C_{11}H_{15}NO_2$	12407-86-2	**I** **M**
	2,3,5-trimethylphenyl methylcarbamate InChI=1/C11H15NO2/c1-7-5-8(2)9(3)10(6-7)14-11(13)12-4/h5-6H,1-4H3,(H,12,13)/f/h12H 3,4,5-trimethylphenyl methylcarbamate InChI=1/C11H15NO2/c1-7-5-10(14-11(13)12-4)6-8(2)9(7)3/h5-6H,1-4H3,(H,12,13)/f/h12H			

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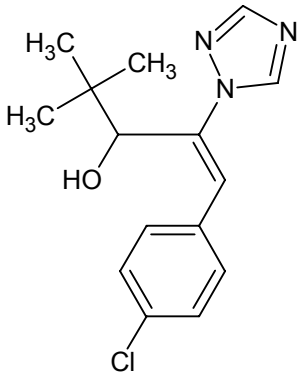
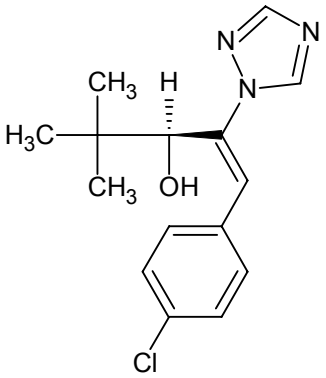
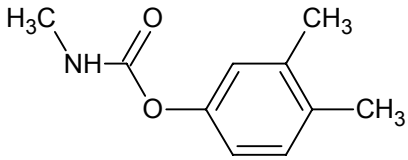
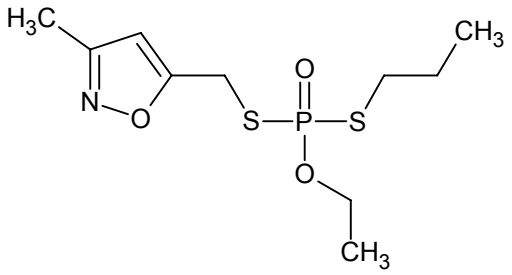
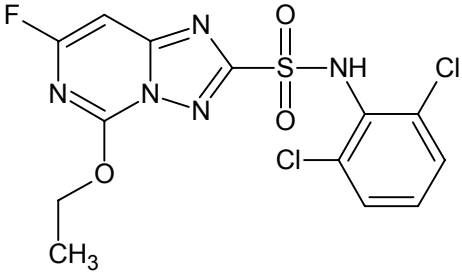
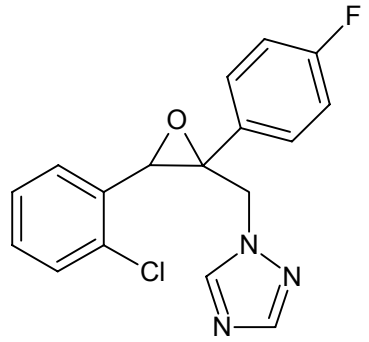
E uniconazole **F** uniconazole (m)	(*E*)-(*RS*)-1-(4-chlorophenyl)-4,4-dimethyl-2-(1*H*-1,2,4-triazol-1-yl)pent-1-en-3-ol (*E*)-(*RS*)-1-(4-chlorophényl)-4,4-diméthyl-2-(1*H*-1,2,4-triazol-1-yl)pent-1-én-3-ol (β*E*)-β-[(4-chlorophenyl)methylene]-α-(1,1-dimethylethyl)-1*H*-1,2,4-triazole-1-ethanol			**F** **P**
		$C_{15}H_{18}ClN_3O$	83657-22-1	
	InChI=1/C15H18ClN3O/c1-15(2,3)14(20)13(19-10-17-9-18-19)8-11-4-6-12(16)7-5-11/h4-10,14,20H,1-3H3/b13-8+/t14-/s3			
E uniconazole-P **F** uniconazole-P (m)	(*E*)-(*S*)-1-(4-chlorophenyl)-4,4-dimethyl-2-(1*H*-1,2,4-triazol-1-yl)pent-1-en-3-ol (*E*)-(*S*)-1-(4-chlorophényl)-4,4-diméthyl-2-(1*H*-1,2,4-triazol-1-yl)pent-1-én-3-ol (α*S*,β*E*)-β-[(4-chlorophenyl)methylene]-α-(1,1-dimethylethyl)-1*H*-1,2,4-triazole-1-ethanol			**F**
		$C_{15}H_{18}ClN_3O$	83657-17-4	
	InChI=1/C15H18ClN3O/c1-15(2,3)14(20)13(19-10-17-9-18-19)8-11-4-6-12(16)7-5-11/h4-10,14,20H,1-3H3/b13-8+/t14-/m1/s1			
E xylylcarb **F** xylylcarb (m)	3,4-xylyl methylcarbamate méthylcarbamate de 3,4-xylyle 3,4-dimethylphenyl methylcarbamate			**I**
		$C_{10}H_{13}NO_2$	2425-10-7	
	InChI=1/C10H13NO2/c1-7-4-5-9(6-8(7)2)13-10(12)11-3/h4-6H,1-3H3,(H,11,12)/f/h11H			

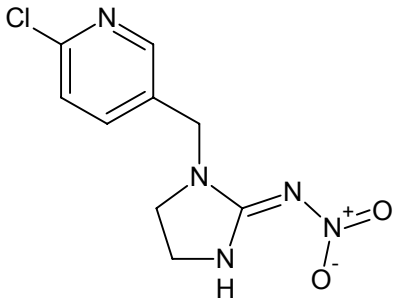
E: Common name F: Nom commun	Chemical name Nom chimique E: IUPAC F: UICPA C: CAS	Structure Structure Molecular formula Formule brute	CAS Registry Number® Numéro d'enregistrement CAS®	Use Applic-ation
	IUPAC International Chemical Identifier (InChI™)			
E zolaprofos **F** zolaprofos (m)	*O*-ethyl *S*-3-methyl-1,2-oxazol-5-ylmethyl *S*-propyl phosphorodithioate phosphorodithioate de *O*-éthyle, de *S*-(3-méthyl-1,2-oxazol-5-yl)méthyle et de *S*-propyle *O*-ethyl *S*-[(3-methyl-5-isoxazolyl)methyl] *S*-propyl phosphorodithioate			I
		$C_{10}H_{18}NO_3PS_2$	63771-69-7	
	InChI=1/C10H18NO3PS2/c1-4-6-16-15(12,13-5-2)17-8-10-7-9(3)11-14-10/h7H,4-6,8H2,1-3H3/t15-/s3			

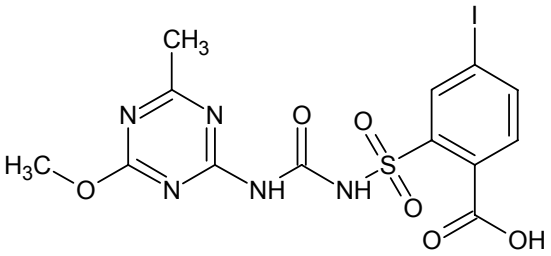
1.2 Corrigenda / Rectificatifs

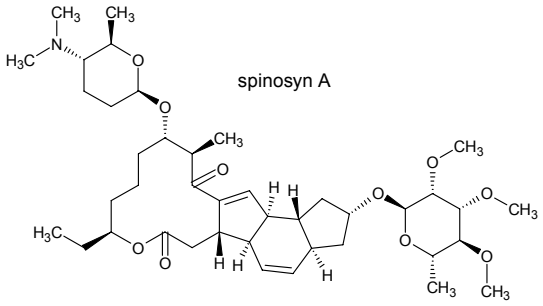
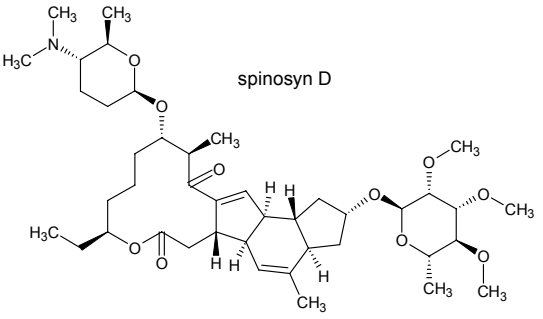
E: Common name F: Nom commun	Chemical name Nom chimique E: IUPAC F: UICPA C: CAS	Structure Structure Molecular formula Formule brute	CAS Registry Number® Numéro d'enregistrement CAS®	Use Applic-ation
	IUPAC International Chemical Identifier (InChI™)			
E diclosulam **F** diclosulame (m)	2′,6′-dichloro-5-ethoxy-7-fluoro[1,2,4]triazolo[1,5-*c*]pyrimidine-2-sulfonanilide *N*-(2,6-dichlorophényl)-5-éthoxy-7-fluoro[1,2,4]triazolo[1,5-*c*]pyrimidine-2-sulfonamide *N*-(2,6-dichlorophenyl)-5-ethoxy-7-fluoro[1,2,4]triazolo[1,5-*c*]pyrimidine-2-sulfonamide			H
		$C_{13}H_{10}Cl_2FN_5O_3S$	145701-21-9	
	InChI=1/C13H10Cl2FN5O3S/c1-2-24-13-17-9(16)6-10-18-12(19-21(10)13)25(22,23)20-11-7(14)4-3-5-8(11)15/h3-6,20H,2H2,1H3			
	NOTE Originally published in Amendment 3 (2001). The previous IUPAC name was not consistent with similar compounds. NOTE Initialement publié dans l'Amendement 3 (2001). Le précédent nom de l'UICPA n'était pas cohérent avec ceux de composés similaires.			

<table>
<tr><th rowspan="2">E: Common name
F: Nom commun</th><th rowspan="2">Chemical name
Nom chimique
E: IUPAC
F: UICPA
C: CAS</th><th colspan="2">Structure
Structure</th><th>Use</th></tr>
<tr><th>Molecular formula
Formule brute</th><th>CAS Registry Number®
Numéro d'enregistrement CAS®</th><th>Applic-ation</th></tr>
<tr><td></td><td colspan="4">IUPAC International Chemical Identifier (InChI™)</td></tr>
<tr><td rowspan="5">E epoxiconazole

F époxiconazole (m)</td><td>(2RS,3SR)-1-[3-(2-chlorophenyl)-2,3-epoxy-2-(4-fluorophenyl)=propyl]-1H-1,2,4-triazole</td><td rowspan="3">  </td><td rowspan="2"></td><td rowspan="5">F</td></tr>
<tr><td>(2RS,3SR)-1-[3-(2-chlorophényl)-2,3-époxy-2-(4-fluorophényl)=propyl]-1H-1,2,4-triazole</td></tr>
<tr><td>rel-1-[[(2R,3S)-3-(2-chloro=phenyl)-2-(4-fluorophenyl)=oxiranyl]methyl]-1H-1,2,4-triazole</td><td>$C_{17}H_{13}ClFN_3O$</td><td>133855-98-8</td></tr>
<tr><td colspan="3">InChI=1/C17H13ClFN3O/c18-15-4-2-1-3-14(15)16-17(23-16,9-22-11-20-10-21-22)12-5-7-13(19)8-6-12/h1-8,10-11,16H,9H2/t16-,17-/s3</td></tr>
<tr><td colspan="3">NOTE Originally published in Amendment 3 (2001). The previous CAS Registry Number was for no stereochemistry.
NOTE Initialement publié dans l'Amendement 3 (2001). Le numéro du CAS précédent ne tenait pas compte de l'aspect stéréochimique.</td></tr>
<tr><td rowspan="5">E imidacloprid

F imidaclopride (m)</td><td>(E)-1-(6-chloro-3-pyridylmethyl)-N-nitroimidazolidin-2-ylideneamine</td><td rowspan="3">  </td><td rowspan="2"></td><td rowspan="5">I</td></tr>
<tr><td>(E)-1-(6-chloro-3-pyridylméthyl)-N-nitroimidazolidin-2-ylidèneamine</td></tr>
<tr><td>(2E)-1-[(6-chloro-3-pyridinyl)methyl]-N-nitro-2-imidazolidinimine</td><td>$C_9H_{10}ClN_5O_2$</td><td>138261-41-3</td></tr>
<tr><td colspan="3">InChI=1/C9H10ClN5O2/c10-8-2-1-7(5-12-8)6-14-4-3-11-9(14)13-15(16)17/h1-2,5H,3-4,6H2,(H,11,13)/f/h11H/b13-9+</td></tr>
<tr><td colspan="3">NOTE Originally published in Amendment 2 (1999). The sponsor has since been able to determine that the product contains almost exclusively the (E)-isomer, and has requested that the definition be amended accordingly.
NOTE Initialement publié dans l'Amendement 2 (1999). Le déclarant a depuis été en mesure de déterminer que le produit contient presque exclusivement l'isomère E et a demandé à ce que la définition soit modifiée en conséquence.</td></tr>
</table>

E: Common name F: Nom commun	Chemical name Nom chimique E: IUPAC F: UICPA C: CAS	Structure Structure		Use
		Molecular formula Formule brute	CAS Registry Number® Numéro d'enregistrement CAS®	Applic-ation
	IUPAC International Chemical Identifier (InChI™)			
E iodosulfuron **F** iodosulfuron (m)	4-iodo-2-(4-methoxy-6-methyl-1,3,5-triazin-2-ylcarbamoyl= sulfamoyl)benzoic acid			**H**
	acide 4-iodo-2-(4-méthoxy-6-méthyl-1,3,5-triazin-2-yl= carbamoylsulfamoyl)benzoïque			
	4-iodo-2-[[[[(4-methoxy-6-methyl-1,3,5-triazin-2-yl)amino]carbonyl]= amino]sulfonyl]benzoic acid	$C_{13}H_{12}IN_5O_6S$	185119-76-0	
	InChI=1/C13H12IN5O6S/c1-6-15-11(18-13(16-6)25-2)17-12(22)19-26(23,24)9-5-7(14)3-4-8(9)10(20)21/h3-5H,1-2H3,(H,20,21)(H2,15,16,17,18,19,22)/f/h17,19-20H			
	NOTE 1 Originally published in Amendment 3 (2001). The previous IUPAC and UICPA names were not consistent with similar compounds. NOTE 2 It should be stated which ester or salt is present, for example iodosulfuron-methyl-sodium [144550-36-7]. NOTE 1 Initialement publié dans l'Amendement 3 (2001). Les noms précédents de l'UICPA et de l'IUPAC n'étaient pas cohérents avec ceux de composés similaires. NOTE 2 Il convient de préciser quel est l'ester ou le sel présent, par exemple iodosulfuron-méthyl-sodium [144550-36-7].			

E: Common name F: Nom commun	Chemical name Nom chimique E: IUPAC F: UICPA C: CAS	Structure Structure		Use
		Molecular formula Formule brute	CAS Registry Number® Numéro d'enregistrement CAS®	Applic-ation
	IUPAC International Chemical Identifier (InChI™)			
E spinosad **F** spinosad (m)	mixture of 50 % to 95 % (2*R*,3a*S*,5a*R*,5b*S*,9*S*,13*S*,14*R*,= 16a*S*,16b*R*)-2-(6-deoxy-2,3,4-tri-O-methyl-α-L-mannopyranosyloxy)-13-(4-dimethylamino-2,3,4,6-tetradeoxy-β-D-erythropyranosyloxy)-9-ethyl-2,3,3a,5a,5b,6,7,9,10,11,12,13,= 14,15,16a,16b-hexadecahydro-14-methyl-1*H*-*as*-indaceno[3,2-*d*]oxacyclododecine-7,15-dione and 50 % to 5 % (2*S*,3a*R*,5a*S*,5b*S*,9*S*,13*S*,14*R*,= 16a*S*,16b*S*)-2-(6-deoxy-2,3,4-tri-O-methyl-α-L-mannopyranosyloxy)-13-(4-dimethylamino-2,3,4,6-tetradeoxy-β-D-erythropyranosyloxy)-9-ethyl-2,3,3a,5a,5b,6,7,9,10,11,12,13,= 14,15,16a,16b-hexadecahydro-4,14-dimethyl-1*H*-*as*-indaceno[3,2-*d*]oxacyclododecine-7,15-dione	 spinosyn A  spinosyn D		I
	mélange constitué à 50–95% de:- (2*R*,3a*S*,5a*R*,5b*S*,9*S*,13*S*,14*R*,= 16a*S*,16b*R*)-2-(6-désoxy-2,3,4-tri-O-méthyl-α-L-mannopyranosyloxy)-13-(4-diméthylamino-2,3,4,6-tétradésoxy-β-D-érythropyranosyloxy)-9-éthyl-2,3,3a,5a,5b,6,7,9,10,11,12,13,= 14,15,16a,16b-hexadécahydro-14-méthyl-1*H*-8-oxacyclo= dodéca[*b*]*as*-indacène-7,15-dione et à 50–5% de:- (2*S*,3a*R*,5a*S*,5b*S*,9*S*,13*S*,14*R*,= 16a*S*,16b*S*)-2-(6-désoxy-2,3,4-tri-O-méthyl-α-L-mannopyranosyloxy)-13-(4-diméthylamino-2,3,4,6-tétradésoxy-β-D-érythropyranosyloxy)-9-éthyl-2,3,3a,5a,5b,6,7,9,10,11,12,13,= 14,15,16a,16b-hexadécahydro-4,14-diméthyl-1*H*-8-oxacyclo= dodéca[*b*]*as*-indacene-7,15-dione			
	(2*R*,3a*S*,5a*R*,5b*S*,9*S*,13*S*,14*R*,= 16a*S*,16b*R*)-2-[(6-deoxy-2,3,4-tri-*O*-methyl-α-L-mannopyranosyl)= oxy]-13-[[(2*R*,5*S*,6*R*)-5-(dimethylamino)tetrahydro-6-methyl-2*H*-pyran-2-yl]oxy]-9-ethyl-2,3,3a,5a,5b,6,9,10,11,12,13,14,= 16a,16b-tetradecahydro-14-methyl-1*H*-*as*-indaceno[3,2-*d*]oxacyclododecin-7,15-dione mixture with (2*S*,3a*R*,5a*S*,5b*S*,9*S*,13*S*,14*R*,= 16a*S*,16b*S*)-2-[(6-deoxy-2,3,4-tri-*O*-methyl-α-L-mannopyranosyl)= oxy]-13-[[(2*R*,5*S*,6*R*)-5-(dimethylamino)tetrahydro-6-methyl-2*H*-pyran-2-yl]oxy]-9-ethyl-2,3,3a,5a,5b,6,9,10,11,12,13,14,= 16a,16b-tetradecahydro-4,14-dimethyl-1*H*-*as*-indaceno[3,2-*d*]oxacyclododecin-7,15-dione	$C_{41}H_{65}NO_{10}$ (spinosyn A) + $C_{42}H_{67}NO_{10}$ (spinosyn D)	168316-95-8	

E: Common name F: Nom commun	Chemical name Nom chimique E: IUPAC F: UICPA C: CAS	Structure Structure		Use
		Molecular formula Formule brute	CAS Registry Number® Numéro d'enregistrement CAS®	Applic-ation
	IUPAC International Chemical Identifier (InChI™)			
	spinosyn A InChI=1/C41H65NO10/c1-10-26-12-11-13-34(52-36-17-16-33(42(5)6)23(3)48-36)22(2)37(44)32-20-30-28(31(32)21-35(43)50-26)15-14-25-18-27(19-29(25)30)51-41-40(47-9)39(46-8)38(45-7)24(4)49-41/h14-15,20,22-31,33-34,36,38-41H,10-13,16-19,21H2,1-9H3/t22-,23-,24+,25-,26+,27-,28-,29-,30-,31+,33+,34+,36+,38+,39-,40-,41+/m1/s1 spinosyn D (4-methyl) InChI=1/C42H67NO10/c1-11-26-13-12-14-35(53-37-16-15-34(43(6)7)24(4)49-37)23(3)38(45)33-20-31-29(32(33)21-36(44)51-26)17-22(2)28-18-27(19-30(28)31)52-42-41(48-10)40(47-9)39(46-8)25(5)50-42/h17,20,23-32,34-35,37,39-42H,11-16,18-19,21H2,1-10H3/t23-,24-,25+,26+,27-,28+,29-,30-,31-,32+,34+,35+,37+,39+,40-,41-,42+/m1/s1		
	NOTE Originally published in Amendment 3 (2001). An error in the IUPAC name is corrected above. NOTE Initialement publié dans l'Amendement 3 (2001). Une erreur dans le nom de l'UICPA est corrigée ci-dessus.			

Index of molecular formulae

Index de formules brutes

$C_{13}H_{15}NO_6$ ..... fenitropan
$C_{13}H_{15}N_3O_3$ ..... imazapyr
$C_{13}H_{16}Cl_3NO_3$ ..... trichlamide
$C_{13}H_{16}NO_4PS$ ..... isoxathion
$C_{13}H_{18}N_2O_2$ ..... hyquincarb
$C_{13}H_{19}ClNO_3PS_2$ ..... anilofos
$C_{13}H_{19}ClNO_3PS_2$ ..... fosmethilan
$C_{13}H_{19}NO$ ..... heptopargil
$C_{13}H_{19}NO_2S$ ..... fenothiocarb
$C_{13}H_{21}O_3PS$ ..... iprobenfos
$C_{13}H_{21}O_4PS$ ..... propaphos
$C_{14}H_5Cl_6NO_3$ ..... tecloftalam
$C_{14}H_6Cl_2F_4N_2O_2$ ..... teflubenzuron
$C_{14}H_7Br_3F_3N_3O_4$ ..... bromethalin
$C_{14}H_8Cl_2N_4$ ..... clofentezine
$C_{14}H_9ClF_4N_2O_4$ ..... flufenpyr
$C_{14}H_9Cl_2NO_5$ ..... bifenox
$C_{14}H_9Cl_5$ ..... *pp'*-DDT
$C_{14}H_{11}Cl_2NO_4$ ..... chlorazifop
$C_{14}H_{12}Cl_2N_2O$ ..... pyrifenox
$C_{14}H_{13}F_3N_6O_5S$ ..... pyroxsulam
$C_{14}H_{14}ClNO_3$ ..... cyprofuram
$C_{14}H_{14}Cl_2N_2O_3$ ..... isovaledione
$C_{14}H_{14}N_4O_5S$ ..... sulfometuron
$C_{14}H_{15}Cl_2N_3OS$ ..... prothioconazole
$C_{14}H_{15}Cl_2N_3O_2$ ..... etaconazole
$C_{14}H_{15}N_5O_6S$ ..... metsulfuron-methyl
$C_{14}H_{16}ClNO_3$ ..... ofurace
$C_{14}H_{16}ClN_3O$ ..... metazachlor
$C_{14}H_{16}ClN_5O_5S$ ..... triasulfuron
$C_{14}H_{17}Cl_2N_3O$ ..... hexaconazole
$C_{14}H_{17}IN_2O_2$ ..... proquinazid
$C_{14}H_{18}ClN_2O_3PS$ ..... pyraclofos
$C_{14}H_{18}N_2O_4$ ..... oxadixyl
$C_{14}H_{19}ClO_3$ ..... mecoprop-P-isobutyl
$C_{14}H_{19}Cl_2FN_2O_3$ ..... fluroxypyr-butometyl
$C_{14}H_{21}NO_3$ ..... furmecyclox
$C_{14}H_{21}NO_4$ ..... diethofencarb
$C_{15}H_8Cl_2F_6N_2O$ ..... flucofuron
$C_{15}H_9ClF_3N_2NaO_6S$ ..... fomesafen-sodium
$C_{15}H_{10}ClF_3N_2O_3$ ..... triflumuron
$C_{15}H_{10}ClF_3N_2O_6S$ ..... fomesafen
$C_{15}H_{11}ClF_3NO_4$ ..... haloxyfop-P
$C_{15}H_{11}Cl_2F_5O_2$ ..... fenfluthrin
$C_{15}H_{12}F_3NO_4$ ..... fluazifop
$C_{15}H_{12}F_3NO_4$ ..... fluazifop-P
$C_{15}H_{14}ClF_4N_3O$ ..... flufenerim
$C_{15}H_{14}Cl_2F_3N_3O_2$ ..... furconazole
$C_{15}H_{15}ClF_3N_3O$ ..... triflumizole
$C_{15}H_{15}ClN_4O_6S$ ..... chlorimuron-ethyl
$C_{15}H_{15}F_3N_2O_2$ ..... flurprimidol
$C_{15}H_{16}N_4O_5S$ ..... sulfometuron-methyl
$C_{15}H_{16}N_4O_7S$ ..... bensulfuron
$C_{15}H_{17}ClN_4$ ..... myclobutanil
$C_{15}H_{17}Cl_2N_3O$ ..... diniconazole
$C_{15}H_{17}Cl_2N_3O$ ..... diniconazole-M
$C_{15}H_{17}Cl_2N_3O_2$ ..... propiconazole
$C_{15}H_{17}N_4NaO_7S$ ..... propoxycarbazone-sodium
$C_{15}H_{18}ClN_3O$ ..... uniconazole
$C_{15}H_{18}ClN_3O$ ..... uniconazole-P
$C_{15}H_{18}Cl_2N_2O_2$ ..... fenoxanil
$C_{15}H_{18}FN_3O_3S$ ..... benthiavalicarb
$C_{15}H_{18}N_4O_7S$ ..... propoxycarbazone
$C_{15}H_{19}Cl_2N_3O$ ..... diclobutrazol
$C_{15}H_{19}N_3O_3$ ..... imazethapyr
$C_{15}H_{20}ClN_3O$ ..... paclobutrazol
$C_{15}H_{21}Cl_2FN_2O_3$ ..... fluroxypyr-meptyl
$C_{15}H_{21}NOS$ ..... dimepiperate
$C_{15}H_{21}NO_4$ ..... metalaxyl-M
$C_{15}H_{22}BrNO$ ..... bromobutide
$C_{15}H_{22}N_4O_3$ ..... imazethapyr-ammonium
$C_{15}H_{23}NO$ ..... tebutam
$C_{15}H_{23}NOS$ ..... esprocarb
$C_{16}H_7ClF_8N_2O_2$ ..... bistrifluron
$C_{16}H_9ClF_3NO_7$ ..... fluoroglycofen
$C_{16}H_{11}ClF_4N_2O_6S$ ..... halosafen
$C_{16}H_{11}ClF_6N_2O$ ..... fluopyram
$C_{16}H_{12}ClF_4N_3O_4$ ..... flumetralin
$C_{16}H_{12}ClNO_4S$ ..... fenthiaprop
$C_{16}H_{12}ClNO_5$ ..... fenoxaprop
$C_{16}H_{12}ClNO_5$ ..... fenoxaprop-P
$C_{16}H_{13}ClFNO_3$ ..... flamprop-M
$C_{16}H_{13}ClF_3NO_4$ ..... haloxyfop-P-methyl
$C_{16}H_{13}F_2N_3O$ ..... flutriafol
$C_{16}H_{14}F_5N_5O_5S$ ..... penoxsulam
$C_{16}H_{14}N_2O_2S$ ..... mefenacet
$C_{16}H_{15}Cl_2NO_2$ ..... clomeprop
$C_{16}H_{15}F_2N_3Si$ ..... flusilazole
$C_{16}H_{18}N_4O_7S$ ..... bensulfuron-methyl
$C_{16}H_{19}N_3$ ..... fenapanil
$C_{16}H_{19}N_5O_9S_2$ ..... mesosulfuron
$C_{16}H_{20}N_2O_2$ ..... benzamacril-isobutyl
$C_{16}H_{23}N_3OS$ ..... buprofezin
$C_{16}H_{24}N_4O_3$ ..... imazapyr-isopropylammonium
$C_{16}H_{25}NOS$ ..... tiocarbazil
$C_{17}H_7Cl_2F_9N_2O_3$ ..... noviflumuron
$C_{17}H_{10}F_6N_4S$ ..... flubenzimine
$C_{17}H_{13}ClFN_3O$ ..... epoxiconazole
$C_{17}H_{13}ClF_3NO_5$ ..... furyloxyfen
$C_{17}H_{13}ClN_2O_4$ ..... quizalofop
$C_{17}H_{13}ClN_2O_4$ ..... quizalofop-P
$C_{17}H_{13}Cl_2NO_4$ ..... chlorazifop-propargyl
$C_{17}H_{14}ClF_7O_2$ ..... tefluthrin
$C_{17}H_{15}ClFNO_3$ ..... flamprop-M-methyl
$C_{17}H_{16}Cl_2N_2O_5$ ..... etnipromid
$C_{17}H_{16}F_3NO_2$ ..... flutolanil
$C_{17}H_{16}O_2$ ..... dofenapyn
$C_{17}H_{17}ClF_4N_4O_5S$ ..... saflufenacil
$C_{17}H_{17}ClN_2O_6S$ ..... fenasulam
$C_{17}H_{17}N_3O_3$ ..... imazaquin
$C_{17}H_{18}F_4O_2$ ..... profluthrin
$C_{17}H_{19}NO_4$ ..... fenoxycarb
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$C_{17}H_{21}ClN_2O_2S$ ..... hexythiazox
$C_{17}H_{21}NO_4S_4$ ..... bensultap
$C_{17}H_{21}N_5O_9S_2$ ..... mesosulfuron-methyl
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$C_{17}H_{24}O_3$ ..... terallethrin
$C_{17}H_{25}NO_5$ ..... alloxydim

$C_{17}H_{26}ClNO_2$ ..... butachlor
$C_{17}H_{27}NO_3S$ ..... cycloxydim
$C_{17}H_{29}NO_3S$ ..... sethoxydim
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$C_{18}H_{12}Cl_2N_2O$ ..... boscalid
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$C_{18}H_{14}Cl_4F_3NO_3$ ..... pyridalyl
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$C_{18}H_{16}ClNO_5$ ..... fenoxaprop-ethyl
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$C_{18}H_{24}N_2O_6$ ..... meptyldinocap
$C_{18}H_{25}N_5O_5$ ..... orysastrobin
$C_{18}H_{26}Cl_2O_3$ 2,4-DB-isoctyl
$C_{18}H_{26}N_2O_5S$ ..... furathiocarb
$C_{18}H_{26}O_2$ ..... cinmethylin
$C_{18}H_{26}O_2$ ..... empenthrin
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$C_{18}H_{41}N_7$ ..... iminoctadine
$C_{19}H_{11}Cl_4N_2NaO_5S$ ..... sulcofuron-sodium
$C_{19}H_{11}F_5N_2O_2$ ..... diflufenican
$C_{19}H_{12}Cl_4N_2O_5S$ ..... sulcofuron
$C_{19}H_{12}F_4N_2O_2$ ..... picolinafen
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$C_{20}H_{17}F_5N_2O_2$ ..... cyflufenamid
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ICS 65.100.01

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