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# Biologische Bundesanstalt

## für Land- und Forstwirtschaft

Merkblatt Nr. 20

4. Auflage

Januar 1981

Verzeichnis der Wirkstoffe  
in  
zugelassenen Pflanzenbehandlungsmitteln

von  
W. Dobrat



Zu beziehen durch:

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Alle Rechte, auch die der Übersetzung, des Nachdrucks und der fotomechanischen Wiedergabe – auch auszugsweise – sind vorbehalten.

Die Wiedergabe von Gebrauchsnamen, Handelsnamen, Warenbezeichnungen usw. in diesem Verzeichnis berechtigt nicht zu der Annahme, daß solche Namen von jedermann benutzt werden dürfen. Es kann sich um gesetzlich geschützte, eingetragene Warenzeichen handeln, auch wenn sie nicht als solche gekennzeichnet sind.

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Dieses Merkblatt enthält eine Zusammenstellung der Wirkstoffe der von der Biologischen Bundesanstalt zugelassenen Pflanzenbehandlungsmittel.

Das Verzeichnis ist alphabetisch nach der gebräuchlichsten Kurzbezeichnung geordnet. Diese Kurzbezeichnung stimmt in den meisten Fällen mit den von der International Organization for Standardization (ISO) empfohlenen oder vorgeschlagenen common names überein. In einigen Fällen wurden common names anderer Organisationen (British Standards Institution, Weed Science Society of America u. a.) oder Namen, die nicht frei verwendbar sind, zum Beispiel Zineb, aufgeführt. Für einige Verbindungen, zum Beispiel Cyanamid, sind die chemischen Bezeichnungen so kurz und verständlich, daß sie keinen common name benötigen. Die ISO-Bezeichnungen wurden bevorzugt verwendet, da in dieser Organisation Staaten aus allen Teilen der Welt zusammenarbeiten. Die common names der ISO stellen jedoch nur Empfehlungen dar und brauchen daher nicht unbedingt von allen Mitgliedstaaten angenommen worden zu sein.

In der Spalte 2 des Verzeichnisses sind die Schlüsselzahlen der Biologischen Bundesanstalt für die Wirkstoffe aufgeführt. Diese Zahlen wurden 1963 eingeführt, weil oft zu einer Zeit, da der common name einer Verbindung noch nicht endgültig festgelegt ist, in der Biologischen Bundesanstalt schon umfangreiche und vielfältige Arbeiten mit den entsprechenden Pflanzenbehandlungsmitteln durchgeführt und Unterlagen und Informationen aller Art registriert werden müssen. So werden in der Biologischen Bundesanstalt alle Unterlagen über die Toxikologie, das Rückstandsverhalten, die Rückstandsanalytik sowie die Analytik des Wirkstoffs in seinen Formulierungen unter dieser Nummer erfaßt, und auch die Literaturkartei wird nach diesem Zahlenschlüssel angelegt.

In der Spalte 3 sind die Schlüsselzahlen des „Collaborative International Pesticides Analytical Council“ (CIPAC) angeführt.

Diese weltweite Organisation für Pflanzenbehandlungsmittel-Analytik schloß sich seinerzeit leider nicht dem deutschen System an, sondern baute unter der Federführung Großbritanniens ein eigenes System neu auf. 1968 wurde dieser Schlüssel von der FAO (Food and Agriculture Organization) übernommen.

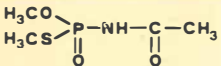
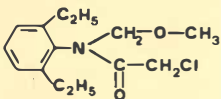
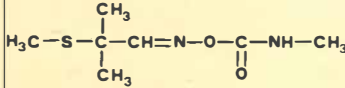
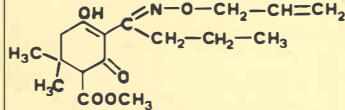

Die Schlüsselzahlen (der Biologischen Bundesanstalt – BBA – bzw. CIPAC) wurden in das Verzeichnis aufgenommen, weil man mit dem BBA-Schlüssel die Analysenmethode für die Rückstände eines Wirkstoffs unmittelbar aus der „Methodensammlung zur Rückstandsanalytik von Pflanzenschutzmitteln“ der Deutschen Forschungsgemeinschaft herausuchen kann, und mit dem CIPAC-Schlüssel die entsprechenden Methoden für die Pflanzenbehandlungsmittelanalyse im „CIPAC Handbook, Analysis of Technical and Formulated Pesticides“ finden kann. Auch die Specifications der FAO sind damit unmittelbar zugänglich.

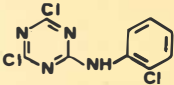
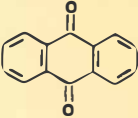
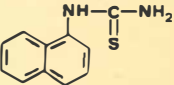
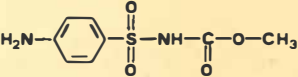
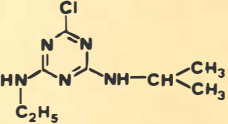
Die chemische Bezeichnung der Wirkstoffe in der Spalte 5 folgt den IUPAC-Regeln. Daneben sind in den nächsten Spalten Strukturformel und Molekulargewicht angegeben.

In der vorletzten Spalte des Verzeichnisses wird der Wirkungsbereich der Verbindung im Pflanzenschutz gekennzeichnet. Hierbei bedeuten die Abkürzungen:

|                 |                                |
|-----------------|--------------------------------|
| A = Akarizid    | O = Avizid                     |
| F = Fungizid    | P = Repellent                  |
| H = Herbizid    | R = Rodentizid                 |
| I = Insektizid  | S = Synergist                  |
| M = Molluskizid | W = Pflanzenwachstumsregulator |
| N = Nematizid   |                                |

In der letzten Spalte werden weitere zur Zeit gebräuchliche Bezeichnungen für den Wirkstoff aufgeführt.

| lfd. Nr. | BBA Nr. | CIPAC Nr. | Name      | Chemische Bezeichnung   | Strukturformel   | Summenformel         | Mol.-gew. | Wir-<br>kungs-<br>bereich | weitere<br>Namen |
|----------|---------|-----------|-----------|---|--|----------------------|-----------|---------------------------|------------------|
| 1        | 358     | 338       | Acephat   | O,S-Dimethyl-N-acetylphosphor-amidothioat                                     |   | $C_4H_{10}NO_3PS$    | 183.2     | I                         |                  |
| 2        | 318     | 204       | Alachlor  | 2'-Chlor-2,6-diethyl-N-methoxymethyl-acetanilid                               |   | $C_{14}H_{20}ClNO_2$ | 269.8     | H                         |                  |
| 3        | 250     | 215       | Aldicarb  | 2-Methyl-2-(methylthio)-propionaldehyd-O-(methylcarbamoyl)oxim                |  | $C_7H_{14}N_2O_2S$   | 190.3     | I<br>N                    |                  |
| 4        | 489     |           | Alloxydim | 2-[(1-Allyloximino)butyl]-4-methoxycarbonyl-5,5-dimethyl-3-oxocyclohex-1-enol |  | $C_{17}H_{25}NO_5$   | 323.4     | H                         |                  |
| 5        | 121     |           |           | Allylalkohol  | $H_2C=CH-CH_2OH$   | $C_3H_6O$            | 58.1      | H                         |                  |
| 6        | 352     | 227       |           | Aluminiumphosphid   | AlP  | AlP                  | 59.0      | I                         |                  |
| 7        | 4       | 90        | Amitrol   | 3-Amino-1,2,4-triazol   |   | $C_2H_4N_4$          | 84.1      | H                         |                  |

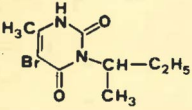
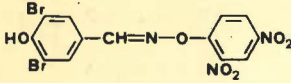
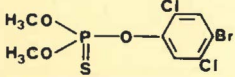
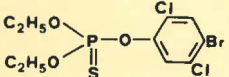

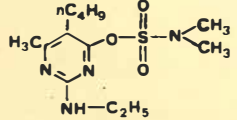
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|----------|---------|-----------|----------|---|--|--------------------|-----------|---------------------------|------------------|
| 7a       | 186     | 294       | Anilazin | 2,4-Dichlor-6-(2-chloranilin)-1,3,5-triazin         |   | $C_9H_5Cl_3N_4$    | 275.5     | F                         | Zinchlor         |
| 8        | 123     | 290       |          | Anthrachinon  |   | $C_{14}H_8O_2$     | 208.2     | P                         |                  |
| 9        | 5       | 39        | Antu     | 1-Naphthylthioharnstoff                             |   | $C_{11}H_{10}N_2S$ | 202.3     | R                         |                  |
| 10       | 312     | 240       | Asulam   | Methyl-N-(4-aminobenzolsulfonyl)-carbamat           |  | $C_8H_{10}O_4N_2S$ | 230.2     | H                         |                  |
| 11       | 6       | 91        | Atrazin  | 4-Ethylamino-6-chlor-2-isopropylamino-1,3,5-triazin |   | $C_8H_{14}ClN_5$   | 215.7     | H                         |                  |

| lfd. Nr. | BBA Nr. | CIPAC Nr. | Name                   | Chemische Bezeichnung  | Strukturformel | Summenformel             | Mol.-gew. | Wir-<br>kungs-<br>bereich | weitere<br>Namen |
|----------|---------|-----------|------------------------|--|----------------|--------------------------|-----------|---------------------------|------------------|
| 12       | 62      | 37<br>b   | Azinphos-ethyl         | O,O-Diethyl-S-(3,4-dihydro-4-oxobenzo-[d]-[1,2,3]-triazin-3-yl)-methyl-dithio-phosphat |                | $C_{12}H_{16}N_3O_3PS_2$ | 345.4     | A<br>I                    |                  |
| 13       | 63      | 37<br>a   | Azinphos-methyl        | S-(3,4-dihydro-4-oxobenzo-[d]-[1,2,3]-triazin-3-yl)-methyl-O,O-dimethyldi-thiophosphat |                | $C_{10}H_{12}N_3O_3PS_2$ | 317.3     | A<br>I                    |                  |
| 14       | 480     |           | Azocyclotin            | Tri(cyclohexyl)-(1,2,4-triazol-1-yl)-zinn  |                | $C_{20}H_{35}N_3Sn$      | 436.2     | A                         |                  |
| 15       | 253     |           | Bacillus Thuringiensis |  |                |                          |           | I                         |                  |
| 16       | 469     | 232       | Bendiocarb             | 2,2-Dimethylbenzo-1,3-dioxolan-4-yl-methylcarbamate                                    |                | $C_{11}H_{13}NO_4$       | 223.2     | I                         |                  |

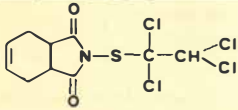
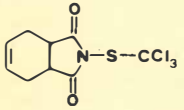
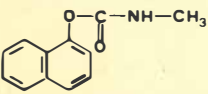
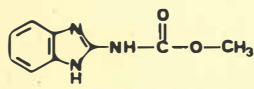
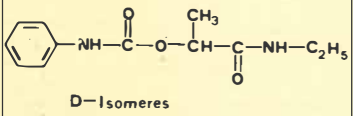
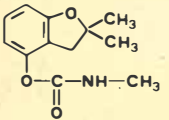


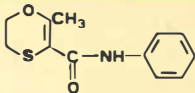
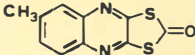
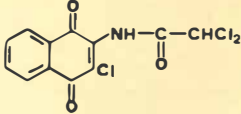
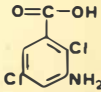
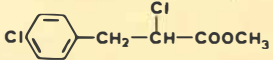
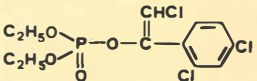
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|----------|---------|-----------|-------------------|---|----------------|------------------------|-----------|---------------------------|------------------|
| 17       | 453     |           | Benodanil         | 2-Jodbenzanilid   |                | $C_{13}H_{10}JNO$      | 323.1     | F                         |                  |
| 18       | 261     | 206       | Benomyl           | Methyl-1-(butyl-carbamoyl)-benzimidazol-2-yl-carbammat  |                | $C_{14}H_{18}N_4O_3$   | 290.3     | F                         |                  |
| 19       | 335     |           | Bentazon          | 3-Isopropyl-(1H)-benzo-2,1,3-thiadiazin-4-on-2,2-dioxid |                | $C_{10}H_{12}N_2O_3S$  | 240.3     | H                         |                  |
| 20       | 360     | 229       | Benzoylprop-ethyl | Ethyl-2-(N-benzoyl-3,4-dichloranilino)-propionat        |                | $C_{18}H_{17}Cl_2NO_3$ | 366.3     | H                         |                  |
| 21       | 8       | 138       | Binapacryl        | 2,4-Dinitro-6-sec. butylphenyl-β-methyl-crotonat        |                | $C_{15}H_{18}N_2O_6$   | 322.4     | A<br>F                    |                  |
| 22       | 11      | 126       | Blausäure         | Cyanwasserstoffsäure                                    | HCN            | CHN                    | 27.0      | I<br>R                    |                  |

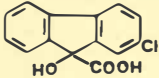
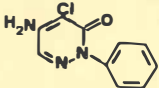
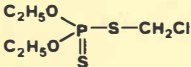
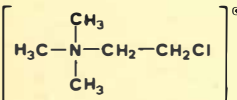
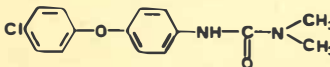
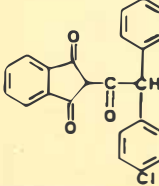


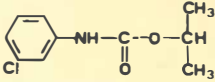
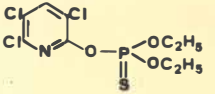

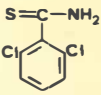
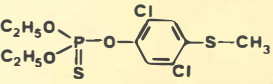
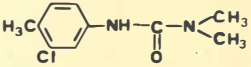
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|----------|---------|-----------|-----------------|---|--|---------------------------|-----------|---------------------------|--------------------|
| 23       | 222     | 139       | Bromacil        | 5-Brom-3-sec-butyl-6-methyl-uracil                              |   | $C_9H_{13}BrN_2O_2$       | 261.1     | H                         |                    |
| 24       | 302     | 269       | Bromfenoxim     | 3,5-Dibrom-4-hydroxybenzaldehyd-2,4-dinitro-phenyloxim          |  | $C_{13}H_7Br_2N_3O_6$     | 461.0     | H                         | Bromofenoxim (ISO) |
| 25       | 210     | 5         | Bromophos       | O-(4-Brom-2,5-dichlor-phenyl)-O,O-dimethyl-thiophosphat         |   | $C_8H_8BrCl_2O_3PS$       | 366.0     | I                         |                    |
| 26       | 263     | 5<br>b    | Bromophos-ethyl | O-(4-Brom-2,5-dichlor-phenyl)-O,O-diethyl-thiophosphat          |   | $C_{10}H_{12}O_3BrCl_2PS$ | 394.0     | A<br>I                    |                    |
| 27       | 264     | 78        | Bromoxynil      | 3,5-Dibrom-4-hydroxy-benzonitril                                |   | $C_7H_3Br_2NO$            | 276.9     | H                         |                    |
| 28       | 418     | 261       | Bupirimat       | 2-Ethylamino-5-n-butyl-6-methylpyrimidin-4-yl-dimethyl-sulfamat |   | $C_{13}H_{24}N_4O_3S$     | 316.4     | F                         |                    |

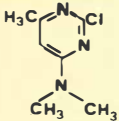
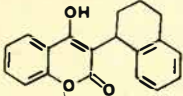
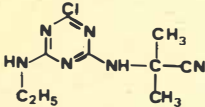
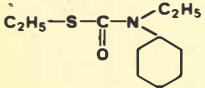
| Ifd. Nr. | BBA Nr. | CIPAC Nr. | Name           | Chemische Bezeichnung                          | Strukturformel  | Summenformel  | Mol.-gew. | Wir-<br>kungs-<br>bereich | weitere<br>Namen |
|----------|---------|-----------|----------------|--|---|---|-----------|---------------------------|------------------|
| 29       | 391     |           | Butocarboxim   | 3-Methylthio-butanon-O-methylcarbamoyloxim     | $\begin{array}{c} \text{CH}_3 \\   \\ \text{H}_3\text{C}-\text{S}-\text{CH}-\text{C}=\text{N}-\text{O}-\overset{\text{O}}{\parallel}{\text{C}}-\text{NH}-\text{CH}_3 \\   \\ \text{CH}_3 \end{array}$   | $\text{C}_7\text{H}_{14}\text{N}_2\text{O}_2\text{S}$ | 190.3     | I                         |                  |
| 30       | 345     |           | Butoxycarboxim | 3-Methylsulfonyl-butanon-O-methylcarbamoyloxim | $\begin{array}{c} \text{O} \quad \text{CH}_3 \\ \parallel \quad   \\ \text{H}_3\text{C}-\text{S}-\text{CH}-\text{C}=\text{N}-\text{O}-\overset{\text{O}}{\parallel}{\text{C}}-\text{NH}-\text{CH}_3 \\ \quad \quad   \\ \quad \quad \text{CH}_3 \end{array}$  | $\text{C}_7\text{H}_{14}\text{N}_2\text{O}_4\text{S}$ | 222.3     | I                         |                  |
| 31       | 361     | 266       | Butylat        | S-Ethyl-diisobutylthiocarbamat                 | $\begin{array}{c} \text{O} \quad \text{CH}_3 \\ \parallel \quad   \\ \text{C}_2\text{H}_5-\text{S}-\text{C}-\text{N}-\text{CH}_2-\text{CH}-\text{CH}_3 \\ \quad \quad \quad \quad   \\ \quad \quad \quad \quad \text{CH}_2-\text{CH}-\text{CH}_3 \\ \quad \quad \quad \quad   \\ \quad \quad \quad \quad \text{CH}_3 \end{array}$ | $\text{C}_{11}\text{H}_{23}\text{NOS}$                | 217.4     | H                         |                  |
| 31a      | 265     |           |                | Calciumcyanamid                                | $\text{Ca}=\text{N}-\text{C}\equiv\text{N}$   | $\text{CaCN}_2$                                       | 80.1      | H<br>F                    | Kalkstickstoff   |
| 32       | 128     | 252       |                | Calciumcyanid                                  | $\begin{array}{c} \text{CN} \\ \diagdown \\ \text{Ca} \\ \diagup \\ \text{CN} \end{array}$  | $\text{Ca}(\text{CN})_2$                              | 92.1      | I                         |                  |
| 33       | 348     |           |                | Calciumphosphid                                | $\begin{array}{c} \text{Ca} \\ \diagdown \quad \diagup \\ \text{P} \\ \diagup \quad \diagdown \\ \text{Ca} \\ \diagdown \quad \diagup \\ \text{P} \\ \diagup \quad \diagdown \\ \text{Ca} \end{array}$  | $\text{Ca}_3\text{P}_2$                               | 182.2     | R                         |                  |

| Ifd. Nr. | BBA Nr. | CIPAC Nr. | Name        | Chemische Bezeichnung  | Strukturformel   | Summenformel         | Mol.-gew. | Wir- kungs- bereich | weitere Namen |
|----------|---------|-----------|-------------|--|--|----------------------|-----------|---------------------|---------------|
| 34       | 266     | 186       | Captafol    | N-(1,1,2,2-Tetra-<br>chloroethyl-thio)-<br>3a,4,7,7a-tetra-<br>hydrophthalimid |   | $C_{10}H_9Cl_4NO_2S$ | 349.1     | F                   |               |
| 35       | 12      | 40        | Captan      | N-(Trichlormethyl-<br>thio)-3a,4,7,7a-<br>tetrahydrophthalimid                 |   | $C_9H_8Cl_3NO_2S$    | 300.6     | F                   |               |
| 36       | 100     | 26        | Carbaryl    | 1-Naphthyl-N-methyl-<br>carbamat   |   | $C_{12}H_{11}NO_2$   | 201.2     | I                   |               |
| 37       | 378     | 263       | Carbendazim | Methyl-benzimidazol-<br>2-yl-carbammat   |  | $C_9H_9N_3O_2$       | 191.2     | F                   | MBC<br>BCM    |
| 38       | 267     | 95        | Carbetamid  | D-1-(Ethylcarbamoyl)-<br>ethylphenylcarbammat                                  |  | $C_{12}H_{16}N_2O_3$ | 236.3     | H                   |               |
|          | 344     | 276       | Carbofuran  | 2,3-Dihydro-2,2-<br>dimethyl-benzo-<br>furan-4-yl-N-methyl-<br>carbammat       |   | $C_{12}H_{15}NO_3$   | 221.3     | I                   |               |

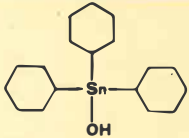
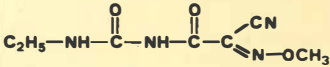
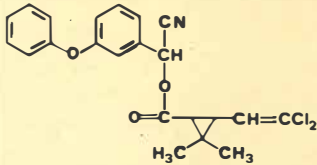
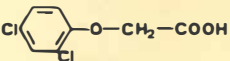
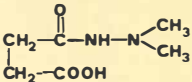
| lfd. Nr. | BBA Nr. | CIPAC Nr. | Name               | Chemische Bezeichnung                               | Strukturformel   | Summenformel           | Mol.-gew. | Wir-<br>kungs-<br>bereich | weitere<br>Namen    |
|----------|---------|-----------|--------------------|---|--|------------------------|-----------|---------------------------|---------------------|
| 40       | 269     | 273       | Carboxin           | 2,4-Dihydro-6-methyl-5-phenylcarbamoyl-1,4-oxathiin |   | $C_{12}H_{13}NO_2S$    | 235.3     | F                         |                     |
| 41       | 398     |           | Cetocetaelat       | Hexadecylpolyglykol-ether                           | $H_3C-(CH_2)_{15}-O-(CH_2-CH_2-O)_8-H$   | $C_{39}H_{66}O_9$      | 594.9     | F                         |                     |
| 42       | 189     | 172       | Chinomethionat     | 6-Methyl-2-oxo-1,3-dithiolo-[4,5-b]-chinoxalin      |   | $C_{10}H_6N_2OS_2$     | 234.3     | F<br>A                    |                     |
| 43       | 442     |           | Chinonamid         | 2,2-Dichlor-N-(3-chlornaphtochinon-2-yl)acetamid    |  | $C_{12}H_6Cl_3NO_3$    | 318.5     | H                         | Quinonamid<br>(ISO) |
| 44       | 342     | 141       | Chloramben         | 3-Amino-2,5-dichlorbenzoessäure                     |   | $C_7H_5Cl_2NO_2$       | 206.0     | H                         |                     |
| 45       | 273     | 270       | Chlorfenpropmethyl | Methyl-2-chlor-3-(4-chlorphenyl)-propionat          |  | $C_{10}H_{10}Cl_2O_2$  | 233.1     | H                         |                     |
| 46       | 239     | 88        | Chlorfenvinphos    | 2-Chlor-1-(2,4 dichlorphenyl)-vinyl-diethylphosphat |  | $C_{12}H_{14}Cl_3O_4P$ | 359.5     | I                         |                     |

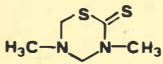
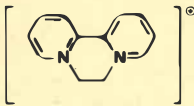
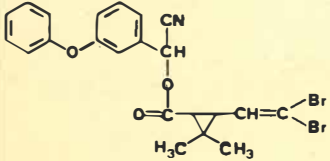
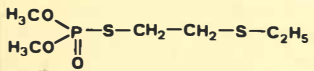
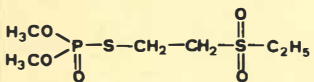
| Ifd. Nr. | BBA Nr. | CIPAC Nr. | Name          | Chemische Bezeichnung   | Strukturformel   | Summenformel           | Mol.-gew. | Wir- kungs- bereich | weitere Namen                |
|----------|---------|-----------|---------------|---|--|------------------------|-----------|---------------------|------------------------------|
| 47       | 275     |           | Chlorflurenol | 2-Chlor-9-hydroxy-<br>fluoren-9-carbonsäure                   |   | $C_{14}H_9ClO_2$       | 274.5     | H<br>W              |                              |
| 48       | 89      | 111       | Chloridazon   | 5-Amino-4-chlor-2-<br>phenylpyridazin-3-on                    |   | $C_{10}H_8ClN_3O$      | 221.6     | H                   | Pyrazon                      |
| 49       | 402     | 337       | Chlormephos   | S-Chlormethyl-O,O-<br>diethyl-dithiophosphat                  |   | $C_5H_{12}ClO_2PS_2$   | 234.6     | I                   |                              |
| 50       | 388     | 143       | Chlormequat   | (2-Chlorethyl)tri-<br>methylammonium                          |   | $C_5H_{13}ClN$         | 158.1     | W                   | CCC<br>(als Chlorid)         |
| 51       | 19      | 187       | Choroxuron    | 3-[4-(4-Chlorphen-<br>oxy)-phenyl]-1,1-di-<br>methylharnstoff |  | $C_{15}H_{15}ClN_2O_2$ | 290.8     | H                   |                              |
| 52       | 238     | 208       | Chlorphacinon | 2-[2-(4-Chlorphenyl)-<br>2-phenyl-acetyl]-<br>indan-1,3-dion  |   | $C_{23}H_{15}ClO_3$    | 374.8     | R                   | Chloro-<br>phacinon<br>(ISO) |

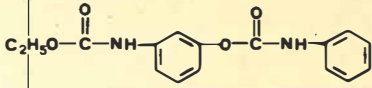
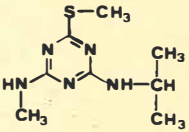
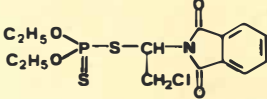
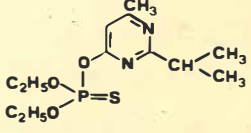
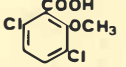

| lfd. Nr. | BBA Nr. | CIPAC Nr. | Name          | Chemische Bezeichnung                                       | Strukturformel   | Summenformel              | Mol.-gew. | Wirkungs-bereich | weitere Namen         |
|----------|---------|-----------|---------------|---|--|---------------------------|-----------|------------------|-----------------------|
| 53       | 21      | 43        | Chlorpropham  | Isopropyl-N-3-chlor-phenylcarbamat                          |   | $C_{10}H_{12}ClNO_2$      | 213.7     | H<br>W           | CIPC                  |
| 54       | 363     | 221       | Chlorpyrifos  | O,O-Diethyl-O-3,5,6-trichlor-2-pyridylthio-phosphat         |   | $C_9H_{11}Cl_3NO_3PS$     | 350.6     | I                |                       |
| 55       | 276     | 288       | Chlorthalonil | Tetrachlorisophthalonitril                                  |   | $C_8Cl_4N_2$              | 265.9     | F                | Chloro-thalonil (ISO) |
| 56       | 224     | 72        | Chlorthiamid  | 2,6-Dichlorthiobenzamid                                     |   | $C_7H_5Cl_2NS$            | 206.1     | H                |                       |
| 57       | 465     |           | Chlorthiophos | O,O-Diethyl-O-2,5-dichlor-4-(methylthio)-phenylthiophosphat |  | $C_{11}H_{15}Cl_2O_3PS_2$ | 360.1     | I                |                       |
| 58       | 279     | 217       | Chlortoluron  | 3-(3-Chlor-4-methyl-phenyl)-1,1-dimethylharnstoff           |  | $C_{10}H_{13}ClN_2O$      | 212.7     | H                |                       |

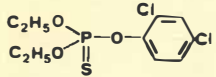
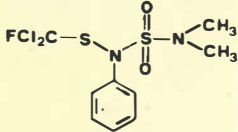
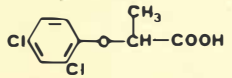
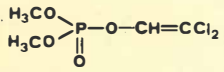
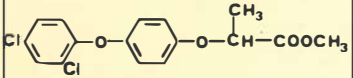
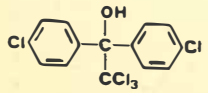
| Ifd. Nr. | BBA Nr. | CIPAC Nr | Name         | Chemische Bezeichnung  | Strukturformel  | Summenformel      | Mol-gew. | Wir-kungs-bereich | weitere Namen        |
|----------|---------|----------|--------------|--|---|-------------------|----------|-------------------|----------------------|
| 59       | 164     | 190      | Crimidin     | 2-Chlor-4-(dimethyl-amino)-6-methyl-pyrimidin                      |  | $C_7H_{10}ClN_3$  | 171,6    | R                 |                      |
| 60       | 26      | 189      | Cumatetralyl | 4-Hydroxy-3-(1,2,3,4-tetrahydro-1-naphthyl)-cumarin                |  | $C_{19}H_{16}O_3$ | 292,3    | R                 | Couma-tetralyl (ISO) |
| 61       | 280     |          |              | Cyanamid   | $H_2N-CN$   | $CH_2N_2$         | 42,0     | H                 |                      |
| 62       | 362     | 230      | Cyanazin     | 4-Ethylamino-2-(1-cyano-1-methylethyl-amino)-6-chlor-1,3,5-triazin |  | $C_9H_{13}ClN_6$  | 240,7    | H                 |                      |
| 63       | 336     | 214      | Cycloat      | N-Ethyl-S-ethyl-N-cyclohexyl-thiocarbamat                          |  | $C_{11}H_{21}NOS$ | 215,4    | H                 |                      |



| lfd. Nr. | BBA Nr. | CIPAC Nr. | Name         | Chemische Bezeichnung  | Strukturformel   | Summenformel           | Mol.-gew. | Wirkungsbereich | weitere Namen |
|----------|---------|-----------|--------------|--|--|------------------------|-----------|-----------------|---------------|
| 64       | 326     | 289       | Cyhexatin    | Trichlohexyl-zinnhydroxid  |   | $C_{18}H_{34}OSn$      | 385.2     | A               |               |
| 64a      | 513     |           | Cymoxanil    | 1-[Cyano(methoximino)acetyl]-3-ethylharnstoff  |  | $C_7H_{10}N_4O_3$      | 198       | F               | Curzate       |
| 65       | 498     | 332       | Cypermethrin | (±)α-Cyano-3-phenoxybenzyl-(±)-cis,trans-3-(2,2-dichlorvinyl)-2,2-dimethylcyclopropanocarboxylat |  | $C_{22}H_{19}Cl_2NO_3$ | 416.3     | I               |               |
| 66       | 27      | 1         | 2,4-D        | 2,4-Dichlorphenoxyessigsäure   |   | $C_8H_6Cl_2O_3$        | 221.0     | H               |               |
| 67       | 28      | 52        | Dalapon      | 2,2-Dichlorpropionsäure  | $H_3C-CCl_2-COOH$  | $C_3H_4Cl_2O_2$        | 143.0     | H               |               |
| 68       | 437     | 330       | Daminozid    | Bernsteinsäuremono(2,2-dimethylhydrazid)   |   | $C_6H_{12}N_2O_3$      | 160.0     | W               |               |

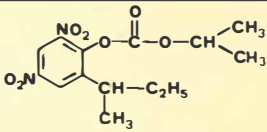
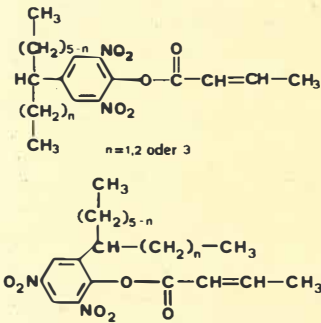
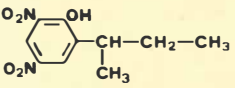
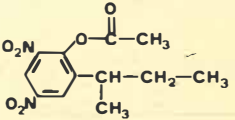
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|----------|---------|------------|------------------------|---|--|------------------------|-----------|-----------------|---------------|
| 69       | 29      | 146        | Dazomet                | Tetrahydro-3,5-dimethyl-1,3,5-thiadiazin-2-thion  |   | $C_5H_{10}N_2S_2$      | 162.3     | N<br>F<br>H     |               |
| 70       | 37      | 55         | Deiquat                | 1,1'-Ethylen-2,2'-bipyridylum   |   | $C_{12}H_{12}N_2$      | 184.2     | H               | Diquat (ISO)  |
| 71       | 496     | 333        | Deltamethrin           | (S)- $\alpha$ -Cyano-3-phenoxybenzyl(1R,3R)-3-(2,2-dibromovinyl)-2,2-dimethylcyclopropanocarboxylat |  | $C_{22}H_{19}Br_2NO_3$ | 505.2     | I               | Decamethrin   |
| 72       | 33      | 47<br>a(S) | Demeton-S-methyl       | S-[2-(Ethylthio)ethyl]-O,O-dimethylthiophosphat   |  | $C_6H_{15}O_3PS_2$     | 230.3     | I<br>A          |               |
| 73       | 77      |            | Demeton-S-methylsulfon | S-[2-(Ethylsulfonyl)ethyl]-O,O-dimethylthiophosphat   |  | $C_6H_{15}O_5PS_2$     | 262.3     | I<br>A          |               |

| lfd. Nr. | BBA Nr. | CIPAC Nr. | Name        | Chemische Bezeichnung  | Strukturformel   | Summenformel             | Mol.-gew. | Wirkungsbereich | weitere Namen |
|----------|---------|-----------|-------------|--|--|--------------------------|-----------|-----------------|---------------|
| 74       | 415     |           | Desmedipham | Ethyl-3-phenyl-carbamoyloxy-phenylcarbamate                      |  | $C_{16}H_{16}N_2O_4$     | 300.3     | H               |               |
| 75       | 244     | 147       | Desmetryn   | 2-Isopropylamino-4-methylamino-6-methylthio-1,3,5-triazin        |   | $C_8H_{15}N_5S$          | 213.3     | H               |               |
| 76       | 281     | 327       | Dialifos    | S-(2-Chlor-1-phthalimidoethyl)-O,O-diethyl-dithiophosphat        |  | $C_{14}H_{17}ClNO_4PS_2$ | 393.8     | I               |               |
| 77       | 35      | 15        | Diazinon    | O,O-Diethyl-O-(2-isopropyl-6-methyl-pyrimidin-4-yl)thio-phosphat |  | $C_{12}H_{21}N_2O_3PS$   | 304.3     | I<br>A          |               |
| 78       | 218     | 85        | Dicamba     | 3,6-Dichlor-2-methoxybenzoësäure                                 |   | $C_8H_6Cl_2O_3$          | 221.0     | H               |               |
| 79       | 225     | 73        | Dichlobenil | 2,6-Dichlorbenzonitril   |   | $C_7H_3Cl_2N$            | 172.0     | H               |               |

| lfd. Nr. | BBA Nr. | CIPAC Nr | Name  | Chemische Bezeichnung   | Strukturformel   | Summenformel  | Mol.-gew. | Wirkungsbereich | weitere Namen |
|----------|---------|----------|---|---|--|---|-----------|-----------------|---------------|
| 80       | 282     | 148      | Dichlofenthion                                | O,O-Diethyl-O-(2,4-dichlorphenyl)-thio-phosphat                                   |   | C <sub>10</sub> H <sub>13</sub> Cl <sub>2</sub> O <sub>3</sub> PS                               | 315.2     | I               |               |
| 81       | 203     | 74       | Dichlofluand                                  | N-[(Dichlorfluor-methyl)-thio]-N',N'-dimethyl-N-phenyl-sulfamid                   |   | C <sub>9</sub> H <sub>11</sub> Cl <sub>2</sub> FN <sub>2</sub><br>O <sub>2</sub> S <sub>2</sub> | 333.2     | F               |               |
| 82       | 38      | 84       | Dichlorprop                                   | 2-(2,4-Dichlorphen-oxyl)propionsäure  |   | C <sub>9</sub> H <sub>8</sub> Cl <sub>2</sub> O <sub>3</sub>                                    | 235.1     | H               | 2,4-DP        |
| 83       | 140     |          |   | 1,3-Dichlorpropen   | ClHC=CH-CH <sub>2</sub> Cl   | C <sub>3</sub> H <sub>4</sub> Cl <sub>2</sub>   | 111.0     | N               |               |
| 84       | 200     | 11       | Dichlorvos                                    | O-2,2-Dichlorvinyl-O,O-dimethylphosphat   |   | C <sub>4</sub> H <sub>7</sub> Cl <sub>2</sub> O <sub>4</sub> P<br>221.0                         | I<br>A    | DDVP            |               |
| 85       | 424     | 358      | Diclofopmethyl                                | Methyl-2-[4-(2,4-dichlorphenoxy)phenoxy]propionat                                 |  | C <sub>16</sub> H <sub>14</sub> Cl <sub>2</sub> O <sub>4</sub>                                  | 341.2     | H               |               |
| 69       | 123     | Dicofol  | 2,2,2-Trichlor-1,1-di-(4-chlorphenyl)-ethanol |  | C <sub>14</sub> H <sub>9</sub> Cl <sub>5</sub> O                                   | 370.5   | A         | Kelthan         |               |

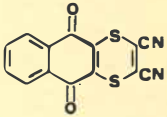
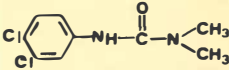
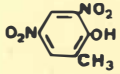
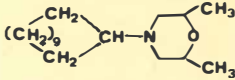
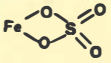
| Ird. Nr. | BBA Nr. | CIPAC Nr. | Name          | Chemische Bezeichnung                                | Strukturformel | Summenformel           | Mol.-gew. | Wirkungsbereich | weitere Namen |
|----------|---------|-----------|---------------|--|----------------|------------------------|-----------|-----------------|---------------|
| 87       | 330     | 299       | Dicrotophos   | 3-(Dimethoxyphosphinyloxy)-N,N-dimethylisocrotonamid |                | $C_8H_{16}NO_5P$       | 237.2     | A<br>I          |               |
| 88       | 285     |           | Dicumarol     | 3,3'-Methylen-bis-(4-hydroxycumarin)                 |                | $C_{19}H_{12}O_6$      | 336.3     | R               |               |
| 89       | 286     |           |               | Dicyclopentadien                                     |                | $C_{10}H_{12}$         | 132.2     | P               |               |
| 90       | 304     |           | Dienochlor    | Perchlorbi-(cyclopenta-2,4-dien-1-yl)                |                | $C_{10}Cl_{10}$        | 474.6     | A               |               |
| 91       | 397     |           | Difenzoquat   | 1,2-Dimethyl-3,5-diphenylpyrazolium                  |                | $C_{17}H_{17}N_2$      | 249.3     | H               |               |
| 92       | 426     | 339       | Diflubenzuron | 1-(4-Chlorphenyl)-3-(2,6-difluorbenzoyl)-harnstoff   |                | $C_{14}H_9ClF_2N_2O_2$ | 310.7     | I               | Difluro       |

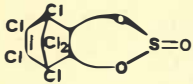
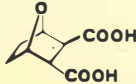
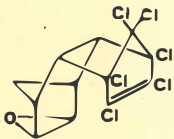
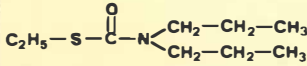
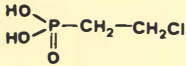
| Ifd. Nr. | BBA Nr. | CIPAC Nr. | Name         | Chemische Bezeichnung  | Strukturformel | Summenformel           | Mol.-gew. | Wirkungsbereich | weitere Namen |
|----------|---------|-----------|--------------|--|----------------|------------------------|-----------|-----------------|---------------|
| 93       | 433     |           | Dikegulac    | 2,3:4,6-Di-O-iso-propyliden- $\alpha$ -L-xylo-2-hexulofuranosonsäure                 |                | $C_{12}H_{18}O_7$      | 274       | W               |               |
| 94       | 40      | 58        | Dimefox      | Bis-(Dimethylamino)-fluorophosphinoxid   |                | $C_4H_{12}FN_2OP$      | 154.1     | A<br>I          |               |
| 95       | 452     | 279       | Dimefuron    | 1-[4-(5-tert. Butyl-2-oxo-1,3,4-oxadiazol-3-yl)-3-chlorphenyl]-3,3-dimethylharnstoff |                | $C_{15}H_{19}ClN_4O_3$ | 338.8     | H               |               |
| 96       | 413     |           | Dimethachlor | $\alpha$ -Chlor-N-(2-methoxyethyl)-2',6'-dimethylacetanilid                          |                | $C_{13}H_{18}ClNO_2$   | 225.7     | H               |               |
| 97       | 42      | 59        | Dimethoat    | O,O-Dimethyl-S-(methylcarbamoylmethyl)-dithiophosphat                                |                | $C_5H_{12}NO_3PS_2$    | 229.3     | A<br>I          |               |
| 98       | 194     |           |              | Dinatriumtetraborat  |                | $Na_2B_4O_7$           | 201.4     | H               | Borax         |

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|----------|---------|-----------|---------------|--|--|----------------------|-----------|------------------|---------------|
| 99       | 255     | 223       | Dinobuton     | 2-sec. Butyl-4,6-dinitrophenyl-isopropylcarbonat   |  | $C_{14}H_{18}N_2O_7$ | 326.3     | A<br>F           |               |
| 100      | 68      | 98        | Dinocap       | Reaktionsgemisch aus 2,6-Dinitro-4-octylphenylcrotonat und 2,4-Dinitro-6-octylphenylcrotonat |  | $C_{18}H_{24}N_2O_6$ | 364.4     | F                |               |
| 101      | 43      | 46        | Dinoseb       | 2-sec. Butyl-4,6-dinitrophenol   |   | $C_{10}H_{12}N_2O_5$ | 240.2     | A<br>F<br>H<br>I |               |
| 102      | 430     |           | Dinosebacetat | 2-sec. Butyl-4,6-dinitrophenylacetat   |   | $C_{12}H_{14}N_2O_6$ | 282.0     | H                |               |

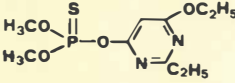
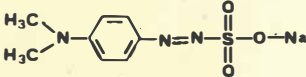
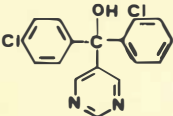
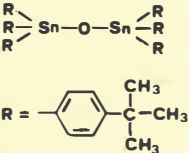
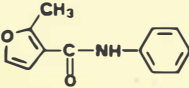


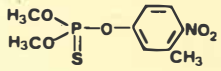
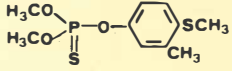
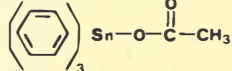
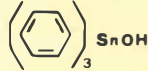
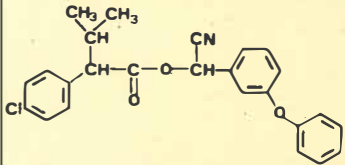
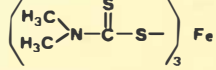
| lfd. Nr. | BBA Nr. | CIPAC Nr. | Name       | Chemische Bezeichnung                                   | Strukturformel | Summenformel            | Mol.-gew. | Wirkungsbereich | weitere Namen |
|----------|---------|-----------|------------|---|----------------|-------------------------|-----------|-----------------|---------------|
| 103      | 333     | 238       | Dinoterb   | 2-tert. Butyl-4,6-dinitrophenol                         |                | $C_{10}H_{12}N_2O_5$    | 240.2     | H               |               |
| 104      | 407     |           | Dioxacarb  | 2-(1,3-Dioxolan-2-yl)-phenyl-methyl-carbamat            |                | $C_{11}H_{13}NO_4$      | 223.2     | I               |               |
| 105      | 31      | 124       | Dioxathion | S,S-1,4-Dioxan-2,3-yliden-bis-(O,O-diethylthiophosphat) |                | $C_{12}H_{26}O_6P_2S_4$ | 456.6     | I               |               |
| 106      | 44      | 152       | Disulfoton | O,O-Diethyl-S-[2-(ethylthio)-ethyl]-dithiophosphat      |                | $C_8H_{19}O_2PS_3$      | 274.4     | A<br>I          |               |
| 107      | 372     |           | Ditalimfos | O,O-Diethylphthalimidothiophosphonat                    |                | $C_{12}H_{14}NO_4PS$    | 299.3     | F               |               |

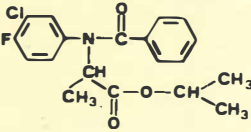
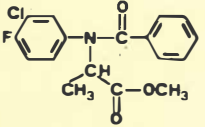
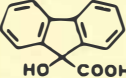
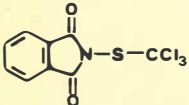
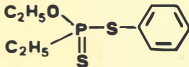
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|----------|---------|-----------|-----------|---|---|----------------------------------|----------------|-----------------|---------------|
| 108      | 45      | 153       | Dithianon | 2,3-Dicyano-1,4-dithia-anthrachinon           |  | $C_{14}H_4N_2O_2S_2$             | 296.3          | F               |               |
| 109      | 46      | 100       | Diuron    | 1-(3,4-Dichlor-phenyl)-3,3-dimethyl-harnstoff |  | $C_9H_{10}Cl_2N_2O$              | 233.1          | H               |               |
| 110      | 47      | 19        | DNOC      | 2-Methyl-4,6-dinitrophenol                    |  | $C_7H_6N_2O_5$                   | 198.1          | H<br>I          |               |
| 111      | 223     | 300       | Dodemorph | 4-Cyclododecyl-2,6-dimethylmorpholin          |  | $C_{18}H_{35}NO$                 | 282.5          | F               |               |
| 112      | 48      | 101       | Dodin     | n-Dodecylguanidinacetat                       | $n-C_{12}H_{25}-NH-C(=NH)-NH_2 \cdot CH_3COOH$                                    | $C_{15}H_{33}N_3O_2$             | 287.5          | F               |               |
| 113      | 229     |           |           | Eisen(II)sulfat                               |  | $FeSO_4$<br>$FeSO_4 \cdot 7H_2O$ | 151.9<br>278.0 | H               |               |

| lfd. Nr. | BBA Nr. | CIPAC Nr. | Name       | Chemische Bezeichnung   | Strukturformel   | Summenformel     | Mol.-gew. | Wirkungsbereich | weitere Namen |
|----------|---------|-----------|------------|---|--|------------------|-----------|-----------------|---------------|
| 114      | 50      | 89        | Endosulfan | 6,7,8,9,10,10-Hexachlor-1,5,5a,6,9,9a-hexahydro-6,9-methano-2,4,3-benzo[e]dioxathiepin-3-oxid     |   | $C_9H_6Cl_6O_3S$ | 407.0     | A<br>I          |               |
| 115      | 395     | 154       | Endothal   | 3,6-Epoxycyclohexan-1,2-dicarbonsäure   |   | $C_8H_{10}O_5$   | 186.1     | H               |               |
| 116      | 52      | 28        | Endrin     | 1,2,3,4,10,10-Hexachlor-6,7-epoxy-1,4,4a,5,6,7,8,8a-octahydro-exo-1,4-exo-5,8-dimethanonaphthalin |   | $C_{12}H_8Cl_6O$ | 380.9     | R<br>(I)        |               |
| 117      | 289     | 155       | EPTC       | S-Ethyl-dipropylthiocarbamat  |  | $C_9H_{19}NOS$   | 189.3     | H               |               |
| 118      | 481     |           | Ethephon   | 2-Chlorethylphosphonsäure   |   | $C_2H_6ClO_3P$   | 144.5     | W               |               |

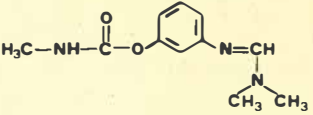
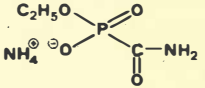
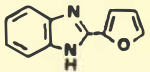
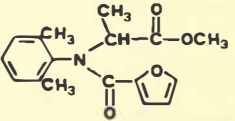
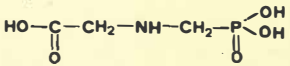
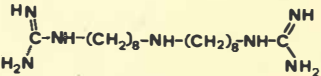
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|----------|---------|-----------|--------------|---|----------------|----------------------|----------|---------------------------|------------------|
| 119      | 406     |           | Ethidimuron  | 1-(5-Ethylsulfonyl-1,3,4-thiadiazol-2-yl)-1,3-dimethylharnstoff |                | $C_7H_{12}N_4O_3S_2$ | 264.3    | H                         | Sulfodiazol      |
| 120      | 393     |           | Ethiofencarb | $\alpha$ -Ethylthio-o-tolyl-methylcarbammat                     |                | $C_{11}H_{15}NO_2S$  | 225.3    | I                         |                  |
| 121      | 290     | 242       | Ethirimol    | 2-Ethylamino-5-n-butyl-4-hydroxy-6-methylpyrimidin              |                | $C_{11}H_{19}N_3O$   | 209.3    | F                         |                  |
| 122      | 383     | 233       | Ethofumesat  | 2-Ethoxy-2,3-dihydro-3,3-dimethylbenzofuran-5-yl-methansulfonat |                | $C_{13}H_{18}O_5S$   | 254.3    | H                         |                  |
| 123      | 387     |           | Etridiazol   | 5-Ethoxy-3-trichloromethyl-1,2,4-thiadiazol                     |                | $C_5H_5Cl_3NOS$      | 247.5    | F                         |                  |

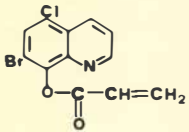
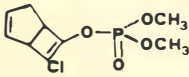
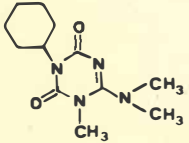
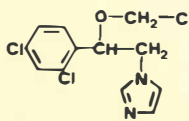
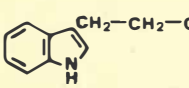
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|----------|---------|-----------|----------------|---|--|--|-----------|---------------------------|------------------|
| 124      | 445     |           | Etrimfos       | O-(6-Ethoxy-2-ethylpyrimidin-4-yl)-O,O-dimethyl-thio-phosphat |   | C <sub>10</sub> H <sub>17</sub> N <sub>2</sub> O <sub>4</sub> PS | 292.0     | I                         |                  |
| 125      | 291     |           | Fenaminosulf   | Natrium-4-dimethyl-aminobenzol-diazo-sulfonat                 |  | C <sub>8</sub> H <sub>10</sub> N <sub>3</sub> NaO <sub>3</sub> S | 251.2     | F                         |                  |
| 126      | 495     |           | Fenarimol      | 2,4'-Dichlor-α-[pyrimidin-5-yl]-benzhydrol                    |   | C <sub>17</sub> H <sub>12</sub> Cl <sub>2</sub> N <sub>2</sub> O | 331.2     | F                         |                  |
| 127      | 410     |           | Fenbutatinoxid | Di[tri-(2-methyl-2-phenylpropyl)-zinn]-oxid                   |   | C <sub>60</sub> H <sub>78</sub> OSn <sub>2</sub>                 | 1052.6    | A                         |                  |
| 128      | 438     |           | Fenfuram       | 2-Methyl-3-furanilid  |   | C <sub>12</sub> H <sub>11</sub> NO <sub>2</sub>                  | 201.0     | F                         |                  |

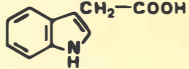
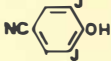
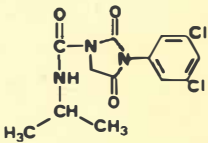
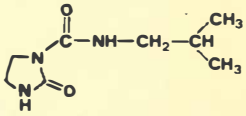
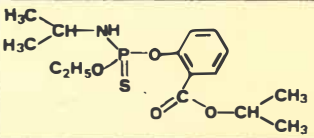
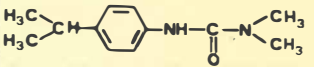
| Ifd. Nr. | BBA Nr. | CIPAC Nr. | Name           | Chemische Bezeichnung                                    | Strukturformel   | Summenformel   | Mol.-gew. | Wirkungsbereich | weitere Namen |
|----------|---------|-----------|----------------|--|--|--|-----------|-----------------|---------------|
| 129      | 58      | 35        | Fenitrothion   | O,O-Dimethyl-O-(4-nitro-m-tolyl)-thiophosphat            |   | C <sub>9</sub> H <sub>12</sub> NO <sub>5</sub> PS              | 277.2     | I<br>A          |               |
| 130      | 57      | 79        | Fenthion       | O,O-Dimethyl-O-(3-methyl-4-methylthiophenyl)thiophosphat |   | C <sub>10</sub> H <sub>15</sub> O <sub>3</sub> PS <sub>2</sub> | 278.3     | I               |               |
| 131      | 55      | 103       | Fentinacetat   | Triphenyl-zinn-acetat                                    |   | C <sub>20</sub> H <sub>18</sub> O <sub>2</sub> Sn              | 409.1     | F               |               |
| 132      | 349     | 103       | Fentinhydroxid | Triphenyl-zinn-hydroxid                                  |   | C <sub>18</sub> H <sub>16</sub> OSn                            | 367.0     | F               |               |
| 133      | 492     | 334       | Fenvalerat     | α-Cyano-3-phenoxybenzyl-isopropyl-p-chlorphenylacetat    |  | C <sub>25</sub> H <sub>22</sub> ClNO <sub>3</sub>              | 419.9     | I               |               |
| 134      | 59      | 57        | Ferbam         | Eisen-tris-(dimethyl-dithiocarbamat)                     |   | C <sub>9</sub> H <sub>18</sub> FeN <sub>3</sub> S <sub>6</sub> | 416.5     | F               |               |

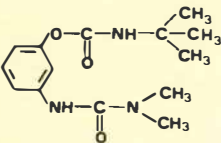
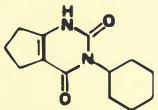

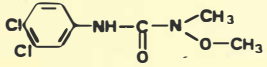
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|----------|---------|-----------|--------------------|---|--|-----------------------|-----------|-----------------|---------------|
| 135      | 468     | 348       | Flamprop-isopropyl | Isopropyl-2-(N-benzoyl-3-chlor-4-fluoranilino)propionat |  | $C_{19}H_{19}ClFNO_3$ | 331.5     | H.              |               |
| 136      | 451     | 349       | Flamprop-methyl    | Methyl-2-(N-benzoyl-3-chlor-4-fluor-anilino)-propionat  |   | $C_{17}H_{15}ClFNO_3$ | 303.5     | H               |               |
| 137      | 215     | 304       | Flurenol           | 9-Hydroxyfluoren-9-carbonsäure                          |   | $C_{14}H_{10}O_3$     | 226.2     | H               |               |
| 138      | 91      | 75        | Folpet             | N-Trichlormethylthio-phthalimid                         |   | $C_9H_4Cl_3NO_2S$     | 296.6     | F               |               |
| 139      | 288     | 306       | Fonofos            | O-Ethyl-S-phenyl-ethyl-dithiophosphonat                 |   | $C_{10}H_{15}OPS_2$   | 246.3     | I               |               |


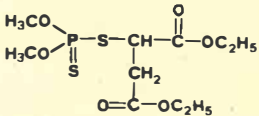
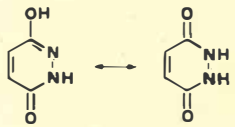
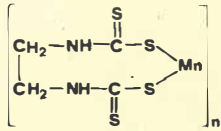


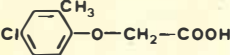
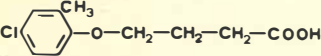
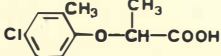
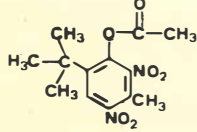
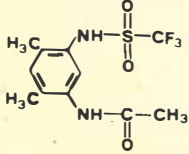
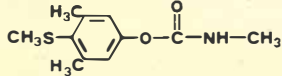
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|----------|---------|-----------|-------------|--|--|----------------------|-----------|-----------------|---------------|
| 140      | 243     |           | Formetanat  | 3-Dimethylamino-methylen-aminophenyl-methyl-carbamat |  | $C_{11}H_{15}N_3O_2$ | 221.3     | A<br>I          |               |
| 141      | 399     | 344       | Fosamin     | Ammonium-ethyl-carbamoylphosphonat                   |   | $C_3H_{11}N_2O_4P$   | 170.1     | H               |               |
| 142      | 214     |           | Fuberidazol | 2-(2-Furyl)-benzimidazol                             |   | $C_{11}H_8N_2O$      | 184.2     | F               |               |
| 142a     | 514     |           | Furalaxyl   | (±)-Methyl-2-[N-(2-furoyl)-N-(2,6-xylyl)]-alaninat   |   | $C_{17}H_{19}NO_4$   | 301.3     | F               |               |
| 143      | 405     | 284       | Glyphosat   | N-(Phosphonomethyl)-glycin                           |  | $C_3H_8NO_5P$        | 169.1     | H               |               |
| 144      | 449     |           | Guazatin    | Di-(8-guanidino-octyl)amin                           |  | $C_{18}H_{41}N_7$    | 355.6     | F               |               |

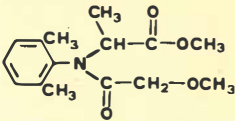
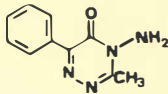
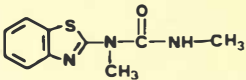
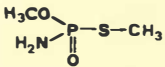
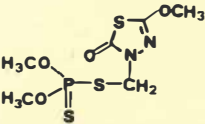
| lfd. Nr. | BBA Nr. | CIPAC Nr. | Name        | Chemische Bezeichnung   | Strukturformel  | Summenformel           | Mol.-gew. | Wirkungsbereich | weitere Namen |
|----------|---------|-----------|-------------|---|---|------------------------|-----------|-----------------|---------------|
| 145      | 420     |           | Halacrinat  | 7-Brom-5-chlor-<br>chinolin-8-yl-acrylat                              |  | $C_{12}H_7BrClNO_2$    | 303.6     | F               |               |
| 146      | 427     |           | Heptenophos | 7-Chlorbicyclo[3,2,0]-<br>hepta-2,6-dien-6-yl-<br>dimethylphosphat    |  | $C_9H_{12}ClO_4P$      | 250.6     |                 |               |
| 147      | 403     |           | Hexazinon   | 3-Cyclohexyl-6-di-<br>methylamino-1-methyl-<br>1,3,5-triazin-2,4-dion |  | $C_{12}H_{20}N_4O_2$   | 252.3     | H               |               |
| 148      | 448     | 335       | Imazalil    | 1-(β-Allyloxy-β-2,4-<br>dichlorphenylethyl)<br>imidazol               |  | $C_{14}H_{14}Cl_2N_2O$ | 297.2     | F               |               |
| 148a     | 145     |           |             | 4-(3-Indol)-buttersäure   |  | $C_{12}H_{13}NO_2$     | 203.2     | W               |               |

| lfd. Nr. | BBA Nr. | CIPAC Nr. | Name        | Chemische Bezeichnung   | Strukturformel   | Summenformel             | Mol.-gew. | Wir-<br>kungs-<br>bereich | weitere<br>Namen |
|----------|---------|-----------|-------------|---|--|--------------------------|-----------|---------------------------|------------------|
| 148b     | 144     |           |             | (3-Indol)-essigsäure  |   | $C_{10}H_9NO_2$          | 175.2     | W                         |                  |
| 149      | 212     | 86        | loxynil     | 4-Hydroxy-3,5-dijodbenzonitril                                      |   | $C_7H_3J_2NO$            | 370.9     | H                         |                  |
| 150      | 419     |           | Iprodion    | 3-(3,5-Dichlorphenyl)-1-isopropylcarbamoyl-imidazolidin-2,4-dion    |   | $C_{13}H_{13}Cl_2N_3O_3$ | 330.2     | F                         | Glycophen        |
| 151      | 359     |           | Isocarbamid | 1-Isobutylcarbamoyl-imidazolidin-2-on                               |   | $C_8H_{15}N_3O_2$        | 185.2     | H                         |                  |
| 152      | 408     |           | Isufenphos  | O-Ethyl-O-(2-isopropoxycarbonyl)phenyl-isopropylphosphoramidothioat |  | $C_{15}H_{24}NO_4PS$     | 345.0     | I                         |                  |
| 153      | 411     | 336       | Isoproturon | 3-(4-Isopropylphenyl)-1,1-dimethylharnstoff                         |  | $C_{12}H_{18}N_2O$       | 206.0     | H                         |                  |

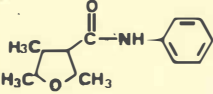
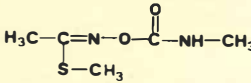
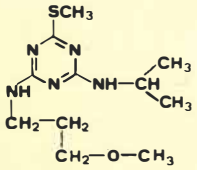
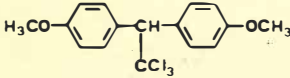
| lfd. Nr. | BBA Nr. | CIPAC Nr. | Name              | Chemische Bezeichnung  | Strukturformel   | Summenformel          | Mol.-gew. | Wir-<br>kungs-<br>bereich | weitere<br>Namen |
|----------|---------|-----------|-------------------|--|--|-----------------------|-----------|---------------------------|------------------|
| 155      | 315     |           | Karbutilat        | 3-(3',3'Dimethyl-ureido)-phenyl-tert. butyl-carbamat             |   | $C_{14}H_{21}N_3O_3$  | 279.3     | H                         |                  |
| 156      | 295     |           | Kupfer-naphthenat |  | Keine definierte Zusammensetzung   |                       |           | F                         |                  |
| 157      | 147     |           |                   | Kupferoxychlorid   | $3 Cu(OH)_2 \cdot CuCl_2$  | $Cu_4H_6Cl_2O_6$      | 427.1     | F                         |                  |
| 158      | 237     | 163       | Lenacil           | 3-Cyclohexyl-6,7-dihydro-1H-cyclopentapyrimidin-2,4-(3H,5H)-dion |   | $C_{13}H_{18}N_2O_2$  | 234.3     | H                         |                  |
| 159      | 70      | 4<br>Y    | Lindan            | $\gamma$ -1,2,3,4,5,6-Hexachlorcyclohexan                        |   | $C_6H_6Cl_6$          | 290.9     | I                         |                  |
| 160      | 71      | 76        | Linuron           | 3-(3,4-Dichlor-phenyl)-1-methoxy-1-methylharnstoff               |  | $C_9H_{10}Cl_2N_2O_2$ | 249.1     | H                         |                  |

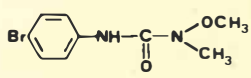
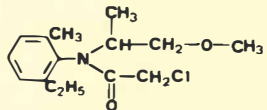
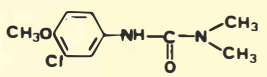
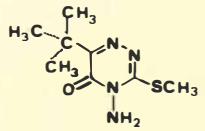
| Ifd. Nr. | BBA Nr. | CIPAC Nr. | Name                | Chemische Bezeichnung  | Strukturformel   | Summenformel          | Mol.-gew. | Wirkungsbereich | weitere Namen |
|----------|---------|-----------|---------------------|--|--|-----------------------|-----------|-----------------|---------------|
| 161      | 354     | 228       |                     | Magnesiumphosphid  |                         | $Mg_3P_2$             | 134.9     | I               |               |
| 162      | 72      | 12        | Malathion           | S-[1,2-bis(Ethoxycarbonyl)-ethyl]-O,O-dimethyldithiophosphat |                        | $C_{10}H_{19}O_6PS_2$ | 330.4     | A<br>I          |               |
| 163      | 297     | 310       | Maleinsäurehydrazid | 1,2-Dihydro-3,6-pyridazinon                                  |                         | $C_4H_4N_2O_2$        | 112.1     | H               |               |
| 164      | 10      | 34        | Mancozeb            | Maneb-Zineb-Mischfällung mit 20 % Mn und 2,5 % Zn            |  |                       |           | F               |               |
| 165      | 73      | 61        | Maneb               | Mangan-ethylen-1,2-bis-dithiocarbamat                        |  <p>Näherungsformel</p> | $C_4H_6MnN_2S_4$      | 265.3     | F               |               |

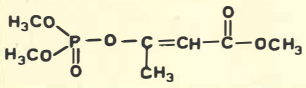

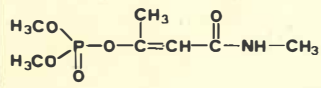
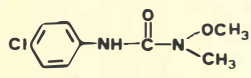
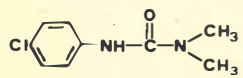
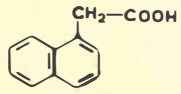
| Ifd. Nr. | BBA Nr. | CIPAC Nr. | Name              | Chemische Bezeichnung  | Strukturformel   | Summenformel   | Mol.-gew. | Wirkungsbereich | weitere Namen |
|----------|---------|-----------|-------------------|--|--|--|-----------|-----------------|---------------|
| 166      | 74      | 2         | MCPA              | 4-Chlor-2-methylphenoxyessigsäure                              |   | C <sub>9</sub> H <sub>9</sub> ClO <sub>3</sub>                                 | 200.6     | H               |               |
| 167      | 75      | 50        | MCPB              | 4-(4-Chlor-2-methylphenoxy)buttersäure                         |  | C <sub>11</sub> H <sub>13</sub> ClO <sub>3</sub>                               | 228.7     | H               |               |
| 168      | 76      | 51        | Mecoprop          | (±)2-(4-Chlor-2-methylphenoxy)propionsäure                     |   | C <sub>10</sub> H <sub>11</sub> ClO <sub>3</sub>                               | 214.6     | H               | MCPP<br>CMPP  |
| 169      | 364     | 164       | Medinoterb-acetat | 2,4-Dinitro-3-methyl-6-tert.butyl-phenyl-acetat                |   | C <sub>13</sub> H <sub>16</sub> N <sub>2</sub> O <sub>6</sub>                  | 296.3     | H               |               |
| 170      | 488     |           | Mefluidid         | N-[2,4-dimethyl-5-(trifluormethylsulfonylamino)phenyl]acetamid |   | C <sub>11</sub> H <sub>13</sub> F <sub>3</sub> N <sub>2</sub> O <sub>3</sub> S | 310.3     | W               |               |
| 79       | 165     |           | Mercapto-dimethur | 4-Methylthio-3,5-xilyl-N-methylcarbamat                        |  | C <sub>11</sub> H <sub>15</sub> NO <sub>2</sub> S                              | 225.3     | I<br>A<br>M     |               |

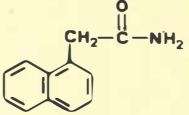
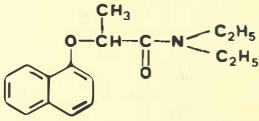
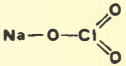
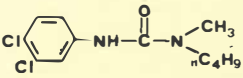
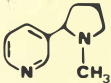
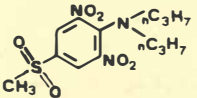
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|----------|---------|-----------|--------------------|---|--|-------------------------|----------------|-----------------|---------------|
| 171a     | 517     |           | Metalaxyl          | Methyl-N-(2-methoxyacetyl)-N-(2,6-xylyl)-alaninat                                       |   | $C_{15}H_{21}NO_4$      | 279.3          | F               | Metaxanin     |
| 172      | 151     | 62        |                    | Metaldehyd  | $(CH_3-CHO)_n$   | $C_8H_{16}O_4$<br>(n=4) | 176.2<br>(n=4) | M               |               |
| 173      | 456     |           | Metamitron         | 4-Amino-4,5-dihydro-3-methyl-6-phenyl-1,2,4-triazin-5-on                                |   | $C_{10}H_{10}N_4O$      | 202.2          | H               |               |
| 174      | 245     | 201       | Methabenzthiazuron | 1,2-Benzothiazolyl-1,3-dimethylharnstoff  |  | $C_{10}H_{11}N_3OS$     | 221.3          | H               |               |
| 175      | 365     | 355       | Methamidophos      | O,S-Dimethylamidothiophosphat   |   | $C_2H_8NO_2PS$          | 141.1          | I               |               |
| 176      | 232     | 193       | Methidathion       | S-(2,3-Dihydro-5-methoxy-2-oxo-1,3,4-thiadiazol-3-ylmethyl)-O,O-dimethyl-dithiophosphat |   | $C_6H_{11}N_2O_4PS_3$   | 302.3          | I               |               |

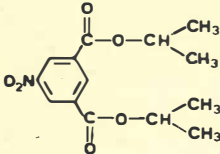
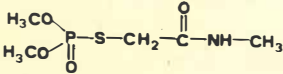
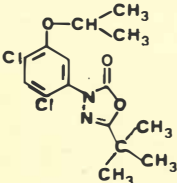
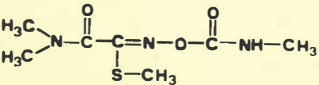


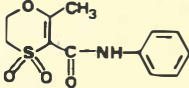
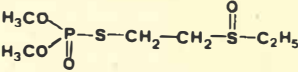
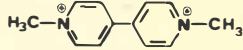
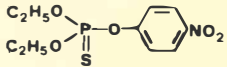
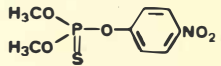
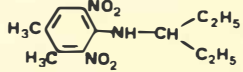
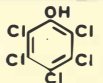
| lfd. Nr. | BBA Nr. | CIPAC Nr. | Name         | Chemische Bezeichnung  | Strukturformel   | Summenformel          | Mol.-gew. | Wirkungsbereich | weitere Namen |
|----------|---------|-----------|--------------|--|--|-----------------------|-----------|-----------------|---------------|
| 176a     | 504     |           | Methfuroxam  | 2,4,5-Trimethyl-3-furanilid  |   | $C_{14}H_{15}NO_2$    | 229.3     | F               | Furavax       |
| 177      | 299     | 264       | Methomyl     | 1-Methylthioethyliden-amino-N-methylcarbamat                             |  | $C_5H_{10}N_2O_2S$    | 162.2     | I               |               |
| 178      | 219     | 94        | Methoprotryn | 2-Isopropylamino-4-(3-Methoxy-n-propyl-amino)-6-methylthio-1,3,5-triazin |   | $C_{11}H_{21}N_5OS$   | 271.4     | H               |               |
| 179      | 80      | 14        | Methoxychlor | 1,1,1-Trichlor-2,2-bis-(4-methoxyphenyl)-ethan                           |  | $C_{16}H_{15}Cl_3O_2$ | 345.7     | I               |               |
| 180      | 149     | 128       | Methylbromid | Brommethan   | $CH_3Br$   | $CH_3Br$              | 94.9      | I<br>N          |               |
|          | 150     |           |              | Methylisothiocyanat  | $H_3C-N=C=S$   | $C_2H_3NS$            | 73.1      | F<br>H<br>N     |               |

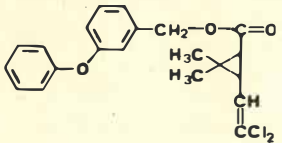
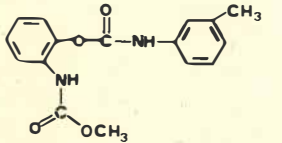
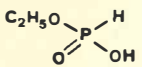
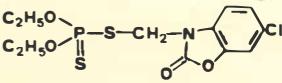
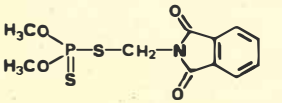
| Ifd. Nr. | BBA Nr. | CIPAC Nr. | Name         | Chemische Bezeichnung  | Strukturformel   | Summenformel  | Mol.-gew. | Wirkungsbereich | weitere Namen |
|----------|---------|-----------|--------------|--|--|---|-----------|-----------------|---------------|
| 182      | 81      |           | Metiram      | Mischfällung bestehend aus dem Ammoniak-Komplex von Zink-(N,N'-ethylenbis(dithiocarbamat) und N,N'-Polyethylenbis-(thiocarbamoyl)-disulfid | $\left[ \begin{array}{c} (-\text{S}-\underset{\text{S}}{\underset{ }{\text{C}}}-\text{NH}-\text{CH}_2-\text{CH}_2-\text{NH}-\underset{\text{S}}{\underset{ }{\text{C}}}-\text{S}-) \cdot \text{Zn}(\text{NH}_3)_3 \\ + \\ (-\text{S}-\underset{\text{S}}{\underset{ }{\text{C}}}-\text{NH}-\text{CH}_2-\text{CH}_2-\text{NH}-\underset{\text{S}}{\underset{ }{\text{C}}}-\text{S}-) \end{array} \right]_n$ | theor. monomere Verbindung<br>$\text{C}_{16}\text{H}_{33}\text{N}_{11}\text{S}_{16}\text{Zn}_3$ | 1088.7    | F               |               |
| 183      | 217     | 168       | Metobromuron | 3-(4-Bromphenyl)-1-methoxy-1-methylharnstoff   |    | $\text{C}_9\text{H}_{11}\text{BrN}_2\text{O}_2$   | 259.1     | H               |               |
| 184      | 422     |           | Metolachlor  | N-(2-Ethyl-6-methylphenyl)-N-(2-methoxy-1-methylethyl)-chloracetamid   |    | $\text{C}_{15}\text{H}_{22}\text{ClNO}_2$   | 283.8     | H               |               |
| 185      | 301     | 219       | Metoxuron    | 3-(3-Chlor-4-methoxyphenyl)-1,1-dimethylharnstoff  |    | $\text{C}_{10}\text{H}_{13}\text{ClN}_2\text{O}_2$  | 228.7     | H               |               |
| 186      | 337     | 283       | Metribuzin   | 4-Amino-4,5-dihydro-3-methylthio-6-tert-butyl-1,2,4-triazin-5-on   |   | $\text{C}_8\text{H}_{14}\text{N}_4\text{OS}$  | 214.3     | H               |               |

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|----------|---------|-----------|---------------|---|--|---------------------|----------|-------------------|---------------|
| 187      | 93      | 45        | Mevinphos     | 2-Methoxycarbonyl-1-methylvinyl-dimethylphosphat  |  | $C_7H_{13}O_6P$     | 224.2    | I<br>A            |               |
| 188      | 191     |           |               | Monochlorbenzol                                   |   | $C_6H_5Cl$          | 112.6    | P                 |               |
| 189      | 259     | 287       | Monocrotophos | 3-Hydroxy-N-methylcrotonamid-O,O-dimethylphosphat |  | $C_4H_{14}NO_5P$    | 223.2    | I<br>A            |               |
| 190      | 82      | 169       | Monolinuron   | 3-(4-Chlorphenyl)-1-methoxy-1-methylharnstoff     |  | $C_9H_{11}ClN_2O_2$ | 214.7    | H                 |               |
| 191      | 83      | 99        | Monuron       | 3-(4-Chlorphenyl)-1,1-dimethylharnstoff           |   | $C_9H_{11}ClN_2O$   | 198.7    | H                 |               |
| 192      | 434     | 313       |               | 1-Naphthylessigsäure                              |   | $C_{12}H_{10}O_2$   | 186.2    | W                 |               |

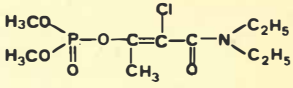
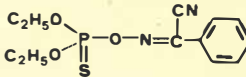
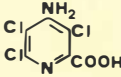
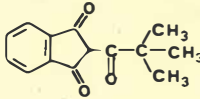
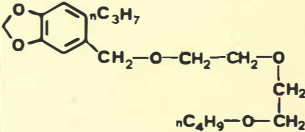
| lfd. Nr. | BBA Nr. | CIPAC Nr. | Name       | Chemische Bezeichnung                             | Strukturformel   | Summenformel           | Mol.-gew. | Wirkungsbereich | weitere Namen |
|----------|---------|-----------|------------|---|--|------------------------|-----------|-----------------|---------------|
| 193      | 435     | 282       |            | 1-Naphthylessigsäure-amid                         |   | $C_{12}H_{11}NO$       | 185.2     | W               |               |
| 194      | 367     | 271       | Napropamid | N,N-Diethyl-2-(1-naphthoxy-propionamid            |  | $C_{17}H_{21}NO_2$     | 271.4     | H               |               |
| 195      | 146     | 7         |            | Natriumchlorat                                    |   | $NaClO_3$              | 106.7     | H               |               |
| 196      | 251     | 314       | Neburon    | 1-n-Butyl-3-(3,4-dichlorphenyl)-1-methylharnstoff |  | $C_{12}H_{16}Cl_2N_2O$ | 275.2     | H               |               |
| 197      | 85      | 8         | Nikotin    | 3-(1-Methylpyrrolidin-2-yl)pyridin                |   | $C_{10}H_{14}N_2$      | 162.2     | I               |               |
| 198      | 368     |           | Nitralin   | 4-Methylsulfonyl-2,6-dinitro-N,N-di-propyl-anilin |   | $C_{13}H_{19}N_3O_6S$  | 345.2     | H               |               |

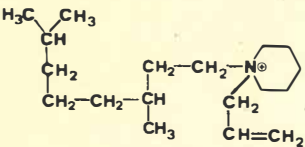
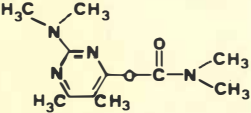
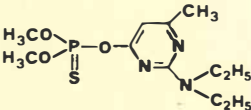
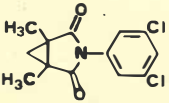
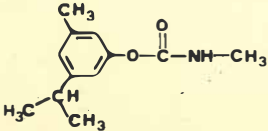
| lfd. Nr. | BBA Nr. | CIPAC Nr. | Name                | Chemische Bezeichnung   | Strukturformel   | Summenformel             | Mol.-gew. | Wirkungsbereich | weitere Namen |
|----------|---------|-----------|---------------------|---|--|--------------------------|-----------|-----------------|---------------|
| 199      | 416     |           | Nitrothaliso-propyl | 5-Nitro-isophthal-säurediisopropylester                                     |   | $C_{14}H_{17}NO_6$       | 295.3     | F               |               |
| 200      | 236     | 202       | Omethoat            | O,O-Dimethyl-S-(N-methylcarbamoyl-methyl)-thiophosphat                      |  | $C_5H_{12}NO_4PS$        | 213.2     | I<br>A          |               |
| 201      | 390     | 213       | Oxadiazon           | 5-tert.-Butyl-3-(2,4-dichlor-5-isopropoxy-phen-1-yl)-1,3,4-oxadiazolin-2-on |   | $C_{15}H_{18}Cl_2N_2O_3$ | 345.2     | H               |               |
| 202      | 441     | 172       | Oxamyl              | 2-Dimethylamino-1-(methylthio)glyoxal-O-methyl-carbamoyl-oxim               |  | $C_7H_{13}N_3O_3S$       | 219.3     | I<br>N          | Thioxamyl     |

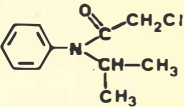
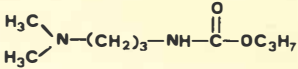
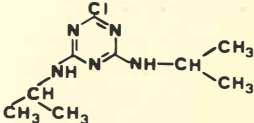
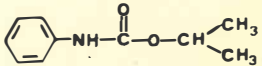
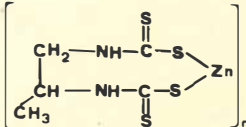
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|----------|---------|-----------|-------------------|--|--|----------------------|-----------|-----------------|--------------------------|
| 203      | 254     | 274       | Oxycarboxin       | 2,3-Dihydro-6-methyl-5-phenylcarbamoyl-1,4-oxathiin-4,4-dioxid |   | $C_{12}H_{13}NO_4S$  | 267.3     | F               |                          |
| 204      | 32      | 171 a     | Oxydemeton-methyl | S-[2-(Ethylsulfinyl)ethyl]-O,O-dimethylthiophosphat            |  | $C_6H_{15}O_4PS_2$   | 246.3     | A               | Demeton-S-methylsulfoxid |
| 205      | 134     | 56        | Paraquat          | 1,1'-Dimethyl-4,4'-bipyridylium                                |  | $C_{12}H_{14}N_2$    | 186.2     | H               |                          |
| 206      | 87      | 10 b      | Parathion         | O,O-Diethyl-O-(4-nitrophenyl)thiophosphat                      |   | $C_{10}H_{14}NO_5PS$ | 291.3     | A<br>I          |                          |
| 207      | 88      | 10 a      | Parathionmethyl   | O,O-Dimethyl-O-(4-nitrophenyl)thiophosphat                     |   | $C_8H_{10}NO_5PS$    | 263.2     | A<br>I          |                          |
| 208      | 404     | 357       | Pendimethalin     | N-(1-Ethylpropyl)-2,6-dinitro-3,4-xylidin                      |  | $C_{13}H_{19}N_3O_4$ | 281.3     | H               | Penoxalin                |
| 209      | 90      | 106       |                   | Pentachlorphenol   |   | $C_6HCl_5O$          | 266.3     | F               |                          |

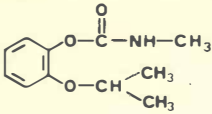
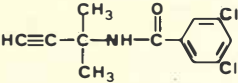
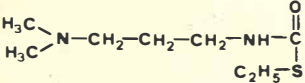
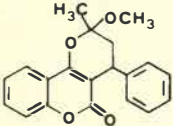
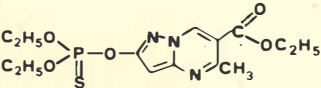
| Ifd. Nr. | BBA Nr. | CIPAC Nr. | Name         | Chemische Bezeichnung   | Strukturformel   | Summenformel             | Mol.-gew. | Wirkungsbereich | weitere Namen |
|----------|---------|-----------|--------------|---|--|--------------------------|-----------|-----------------|---------------|
| 210      | 494     | 331       | Permethrin   | 3-Phenoxybenzyl-(±)-cis-trans-3-(2,2-dichlorvinyl)-2,2-dimethylcyclopropan-carboxylat |  | $C_{21}H_{20}Cl_2O_3$    | 391.3     | I               |               |
| 211      | 233     | 77        | Phenmedipham | Methyl-3-(m-tolyl-carbamoyloxý)-phenylcarbamát  |  | $C_{16}H_{16}N_2O_4$     | 300.3     | H               |               |
| 211a     |         |           | Phosethyl    | Ethylhydrogen-phosphonat  |   | $C_2H_7O_3P$             | 110.0     | F               | Efosit        |
| 212      | 306     | 109       | Phosalon     | S-(6-Chlor-2-oxo-benzoxazolin-3-yl) methyl-O,O-diethyl-dithiophosphat                 |  | $C_{12}H_{15}ClNO_4PS_2$ | 367.8     | A               |               |
| 213      | 334     | 318       | Phosmet      | O,O-Dimethyl-S-phthalimidomethyl-dithiophosphat                                       |  | $C_{11}H_{12}NO_4S_2P$   | 317.3     | A<br>I          |               |

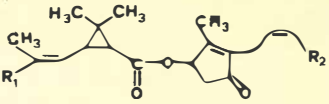
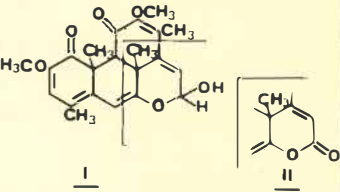


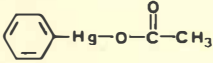
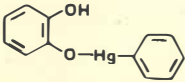
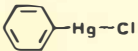

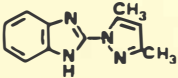
| Ifd. Nr. | BBA Nr. | CIPAC Nr. | Name             | Chemische Bezeichnung   | Strukturformel   | Summenformel           | Mol.-gew. | Wirkungsbereich | weitere Namen |
|----------|---------|-----------|------------------|---|--|------------------------|-----------|-----------------|---------------|
| 214      | 94      | 110       | Phosphamidon     | O-(2-Chlor-2-diethyl-carbamoyl-1-methyl-vinyl)-O,O-dimethylphosphat |  | $C_{10}H_{19}ClNO_5P$  | 299.7     | A<br>I          |               |
| 215      | 307     |           | Phoxim           | O- $\alpha$ -Cyanobenzyliden-amino-O,O-diethylthiophosphat          |  | $C_{12}H_{15}N_2O_3PS$ | 298.3     | I               |               |
| 216      | 308     | 174       | Picloram         | 4-Amino-3,5,6-trichlorpicolinsäure                                  |   | $C_6H_3Cl_3N_2O_2$     | 241.5     | H               |               |
| 217      | 95      | 196       | Pindon           | 2-Pivaloylindan-1,3-dion  |   | $C_{14}H_{14}O_3$      | 230.3     | R               |               |
| 218      | 163     | 33        | Piperonylbutoxid | 5-[2-(2-Butoxyethoxy)-ethoxymethyl]-6-propyl-1,3-benzodioxol        |  | $C_{19}H_{30}O_5$      | 338.5     | S               |               |

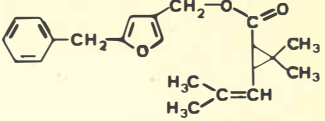
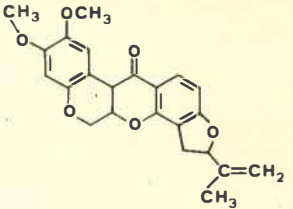
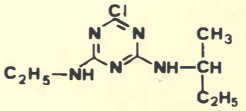
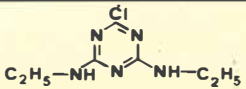
| lfd. Nr. | BBA Nr. | CIPAC Nr. | Name             | Chemische Bezeichnung  | Strukturformel   | Summenformel   | Mol.-gew. | Wirkungsbereich | weitere Namen |
|----------|---------|-----------|------------------|--|--|--|-----------|-----------------|---------------|
| 219      | 432     |           | Piproctanyl      | 1-Allyl-1(3,7-dimethyloctyl)-piperidinium                          |  | C <sub>18</sub> H <sub>36</sub> N                                | 266.4     | W               |               |
| 220      | 309     | 231       | Pirimicarb       | 2-Dimethylamino-5,6-dimethylpyrimidin-4-yl-dimethylcarbamate       |   | C <sub>11</sub> H <sub>18</sub> N <sub>4</sub> O <sub>2</sub>    | 238.3     | I               |               |
| 221      | 476     | 239 a     | Pirimiphosmethyl | O-(2-Diethylamino-6-methylpyrimidin-4-yl)-O,O-dimethylthiophosphat |   | C <sub>11</sub> H <sub>20</sub> N <sub>3</sub> O <sub>3</sub> PS | 305.0     | A<br>I          |               |
| 222      | 491     |           | Procymidon       | N-(3,5-Dichlorphenyl)-1,2-dimethyl-1,2-cyclopropanedicarboximid    |   | C <sub>13</sub> H <sub>11</sub> Cl <sub>2</sub> NO <sub>2</sub>  | 284.1     | F               |               |
| 223      | 190     |           | Promecarb        | 3-Isopropyl-5-methylphenyl-methylcarbamate                         |  | C <sub>12</sub> H <sub>17</sub> NO <sub>2</sub>                  | 207.3     | I               |               |

| Ird. Nr. | BBA Nr. | CIPAC Nr. | Name        | Chemische Bezeichnung                         | Strukturformel   | Summenformel         | Mol.-gew             | Wir-<br>kungs-<br>bereich | weitere<br>Namen |
|----------|---------|-----------|-------------|---|--|----------------------|----------------------|---------------------------|------------------|
| 224      | 310     | 176       | Propachlor  | 2'-Chlor-N-iso-propylacetanilid               |   | $C_{11}H_{14}ClNO$   | 211.7                | H                         |                  |
| 224a     | 516     |           | Propamocarb | Propyl-3-(dimethyl-amino)-propylcarbamate     |  | $C_9H_{21}ClN_2O_2$  | 224.7                | F                         |                  |
| 225      | 97      | 92        | Propazin    | 6-Chlor-2,4-bis-iso-propylamino-1,3,5-triazin |  | $C_9H_{10}ClN_5$     | 229.7                | H                         |                  |
| 226      | 66      | 63        | Propham     | Isopropyl-N-phenyl-carbamate                  |  | $C_{10}H_{13}NO_2$   | 179.2                | H<br>W                    | IPC              |
| 227      | 117     | 177       | Propineb    | Zink-propylen-bis-dithiocarbamat              |  | $(C_5H_8N_2S_4Zn)_n$ | (289.8) <sub>n</sub> | F                         |                  |

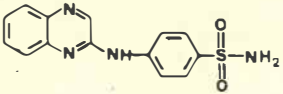
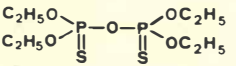
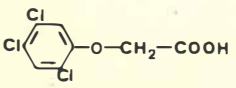
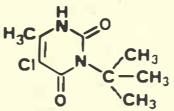
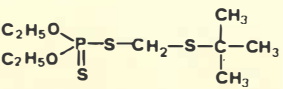
| Ifd. Nr. | BBA Nr. | CIPAC Nr. | Name          | Chemische Bezeichnung  | Strukturformel   | Summenformel   | Mol.-gew. | Wir- kungs- bereich | weitere Namen |
|----------|---------|-----------|---------------|--|--|--|-----------|---------------------|---------------|
| 228      | 216     | 80        | Propoxur      | 2-Isopropoxyphenyl-methylcarbamate   |   | C <sub>11</sub> H <sub>15</sub> NO <sub>3</sub>                  | 209.2     | I                   |               |
| 229      | 350     | 315       | Propyzamid    | 3,5-Dichlor-N-(1,1-dimethylpropinyl)-benzamid  |   | C <sub>12</sub> H <sub>11</sub> Cl <sub>2</sub> NO               | 256.1     | H                   |               |
| 230      | 396     |           | Prothiocarb   | S-Ethyl-N-(3-dimethylaminopropyl)-thiocarbamat   |  | C <sub>8</sub> H <sub>18</sub> N <sub>2</sub> OS                 | 190.3     | F                   |               |
| 231      | 132     |           | Pyranocumarin | 2'-Methoxy-2'-methyl-4'-phenyl-2,3-dihydro-4H,5H-pyrano-[3,2-c]-[1]-benzopyran-5-on    |   | C <sub>20</sub> H <sub>18</sub> O <sub>4</sub>                   | 322.4     | R                   |               |
| 232      | 328     | 350       | Pyrazophos    | O,O-Diethyl-O-(6-ethoxy-carbonyl-5-methyl-pyrazolo-[1,5a]-pyrimidin-2-yl)-thiophosphat |  | C <sub>14</sub> H <sub>20</sub> N <sub>3</sub> O <sub>5</sub> PS | 373.4     | F                   |               |

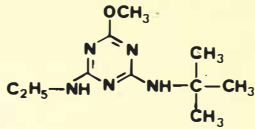
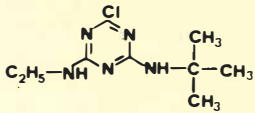
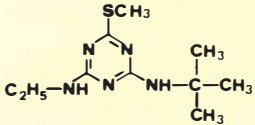
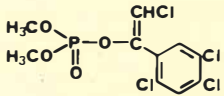
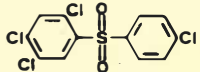
| Ifd. Nr.       | BBA Nr.           | CIPAC Nr.                | Name                                   | Chemische Bezeichnung  | Strukturformel  | Summenformel  | Mol.-gew.        | Wirkungsbereich  | weitere Namen |               |                |                          |  |                |                   |                          |  |             |                |                |  |              |                   |                |  |              |                |                         |  |               |                   |                         |  |  |  |   |  |
|----------------|-------------------|--------------------------|--|--|---|---|------------------|------------------|---------------|---------------|----------------|--------------------------|--|----------------|-------------------|--------------------------|--|-------------|----------------|----------------|--|--------------|-------------------|----------------|--|--------------|----------------|-------------------------|--|---------------|-------------------|-------------------------|--|--|--|---|--|
| 233            | 98                | 32                       | Pyrethrine                             | Gemisch aus Pyrethrin I und II<br>Cinerin I und II<br>Jasmolin I und II                  |  <p>allgemeines Strukturschema</p> <table data-bbox="718 376 1047 540"> <thead> <tr> <th></th> <th><math>\overline{R_1}</math></th> <th><math>\overline{R_2}</math></th> <th></th> </tr> </thead> <tbody> <tr> <td>Pyrethrin I :</td> <td><math>-\text{CH}_3</math></td> <td><math>-\text{CH}=\text{CH}_2</math></td> <td><math>\text{C}_{21}\text{H}_{28}\text{O}_3</math></td> </tr> <tr> <td>Pyrethrin II :</td> <td><math>-\text{COOCH}_3</math></td> <td><math>-\text{CH}=\text{CH}_2</math></td> <td><math>\text{C}_{22}\text{H}_{28}\text{O}_5</math></td> </tr> <tr> <td>Cinerin I :</td> <td><math>-\text{CH}_3</math></td> <td><math>-\text{CH}_3</math></td> <td><math>\text{C}_{20}\text{H}_{28}\text{O}_3</math></td> </tr> <tr> <td>Cinerin II :</td> <td><math>-\text{COOCH}_3</math></td> <td><math>-\text{CH}_3</math></td> <td><math>\text{C}_{21}\text{H}_{28}\text{O}_5</math></td> </tr> <tr> <td>Jasmolin I :</td> <td><math>-\text{CH}_3</math></td> <td><math>-\text{C}_2\text{H}_5</math></td> <td><math>\text{C}_{21}\text{H}_{30}\text{O}_3</math></td> </tr> <tr> <td>Jasmolin II :</td> <td><math>-\text{COOCH}_3</math></td> <td><math>-\text{C}_2\text{H}_5</math></td> <td><math>\text{C}_{22}\text{H}_{30}\text{O}_5</math></td> </tr> </tbody> </table> |   | $\overline{R_1}$ | $\overline{R_2}$ |               | Pyrethrin I : | $-\text{CH}_3$ | $-\text{CH}=\text{CH}_2$ | $\text{C}_{21}\text{H}_{28}\text{O}_3$ | Pyrethrin II : | $-\text{COOCH}_3$ | $-\text{CH}=\text{CH}_2$ | $\text{C}_{22}\text{H}_{28}\text{O}_5$ | Cinerin I : | $-\text{CH}_3$ | $-\text{CH}_3$ | $\text{C}_{20}\text{H}_{28}\text{O}_3$ | Cinerin II : | $-\text{COOCH}_3$ | $-\text{CH}_3$ | $\text{C}_{21}\text{H}_{28}\text{O}_5$ | Jasmolin I : | $-\text{CH}_3$ | $-\text{C}_2\text{H}_5$ | $\text{C}_{21}\text{H}_{30}\text{O}_3$ | Jasmolin II : | $-\text{COOCH}_3$ | $-\text{C}_2\text{H}_5$ | $\text{C}_{22}\text{H}_{30}\text{O}_5$ |  |  | I |  |
|                | $\overline{R_1}$  | $\overline{R_2}$         |  |  |   |   |                  |                  |               |               |                |                          |  |                |                   |                          |  |             |                |                |  |              |                   |                |  |              |                |                         |  |               |                   |                         |  |  |  |   |  |
| Pyrethrin I :  | $-\text{CH}_3$    | $-\text{CH}=\text{CH}_2$ | $\text{C}_{21}\text{H}_{28}\text{O}_3$ |  |   |   |                  |                  |               |               |                |                          |  |                |                   |                          |  |             |                |                |  |              |                   |                |  |              |                |                         |  |               |                   |                         |  |  |  |   |  |
| Pyrethrin II : | $-\text{COOCH}_3$ | $-\text{CH}=\text{CH}_2$ | $\text{C}_{22}\text{H}_{28}\text{O}_5$ |  |   |   |                  |                  |               |               |                |                          |  |                |                   |                          |  |             |                |                |  |              |                   |                |  |              |                |                         |  |               |                   |                         |  |  |  |   |  |
| Cinerin I :    | $-\text{CH}_3$    | $-\text{CH}_3$           | $\text{C}_{20}\text{H}_{28}\text{O}_3$ |  |   |   |                  |                  |               |               |                |                          |  |                |                   |                          |  |             |                |                |  |              |                   |                |  |              |                |                         |  |               |                   |                         |  |  |  |   |  |
| Cinerin II :   | $-\text{COOCH}_3$ | $-\text{CH}_3$           | $\text{C}_{21}\text{H}_{28}\text{O}_5$ |  |   |   |                  |                  |               |               |                |                          |  |                |                   |                          |  |             |                |                |  |              |                   |                |  |              |                |                         |  |               |                   |                         |  |  |  |   |  |
| Jasmolin I :   | $-\text{CH}_3$    | $-\text{C}_2\text{H}_5$  | $\text{C}_{21}\text{H}_{30}\text{O}_3$ |  |   |   |                  |                  |               |               |                |                          |  |                |                   |                          |  |             |                |                |  |              |                   |                |  |              |                |                         |  |               |                   |                         |  |  |  |   |  |
| Jasmolin II :  | $-\text{COOCH}_3$ | $-\text{C}_2\text{H}_5$  | $\text{C}_{22}\text{H}_{30}\text{O}_5$ |  |   |   |                  |                  |               |               |                |                          |  |                |                   |                          |  |             |                |                |  |              |                   |                |  |              |                |                         |  |               |                   |                         |  |  |  |   |  |
| 234            | 258               |                          | Quassin                                | Molekülverbindung<br>1:1 aus Neo-<br>quassin (I) und<br>Isoquassin (II)<br>nach Polonsky |   | $\text{C}_{22}\text{H}_{30}\text{O}_6$ (I)<br>$\text{C}_{22}\text{H}_{28}\text{O}_6$ (II) | 780.9<br>778.9   | I<br>P           |               |               |                |                          |  |                |                   |                          |  |             |                |                |  |              |                   |                |  |              |                |                         |  |               |                   |                         |  |  |  |   |  |
| 235            | 173               |                          |  | 2-Methoxyethyl-Hg-<br>chlorid  | $\text{H}_3\text{C}-\text{O}-\text{CH}_2-\text{CH}_2-\text{Hg}-\text{Cl}$   | $\text{C}_3\text{H}_7\text{ClHgO}$  | 295.1            | F                |               |               |                |                          |  |                |                   |                          |  |             |                |                |  |              |                   |                |  |              |                |                         |  |               |                   |                         |  |  |  |   |  |

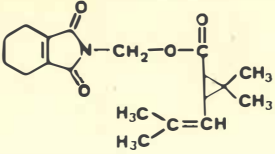
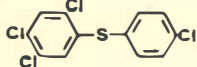
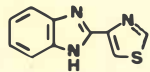
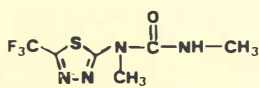
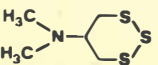
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|----------|---------|-----------|------------|--|---|---|-----------|-----------------|-----------------|
| 236      | 172     |           |            | 2-Methoxyethyl-Hg-silicat                        | $\text{H}_3\text{C}-\text{O}-\text{CH}_2-\text{CH}_2-\text{Hg}-\text{silicat}$    | Genauere Formel nicht bekannt. Das technische Produkt enthält 30 % Hg |           | F               |                 |
| 237      | 180     |           |            | Phenyl-Hg-acetat                                 |  | $\text{C}_8\text{H}_8\text{HgO}_2$                                    | 336.8     | F               |                 |
| 238      | 179     |           |            | Phenyl-Hg-brenz-catechin                         |  | $\text{C}_{12}\text{H}_{10}\text{O}_2\text{Hg}$                       | 386.2     | F               |                 |
| 239      | 181     |           |            | Phenyl-Hg-chlorid                                |  | $\text{C}_6\text{H}_5\text{ClHg}$                                     | 295.3     | F               |                 |
| 240      | 99      | 78        | Quintozen  | Pentachlornitrobenzol                            |  | $\text{C}_6\text{Cl}_5\text{NO}_2$                                    | 295.4     | F               |                 |
| 240a     | 606     |           | Rabenzazol | 2-(3,5-Dimethyl-1H-pyrazol-1-yl)-1H-benzimidazol |  | $\text{C}_{12}\text{H}_{12}\text{N}_4$                                | 213.3     | F               | Dimeth-benzazol |

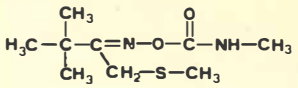
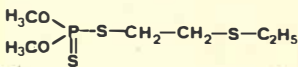
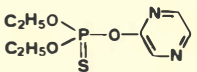
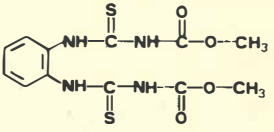
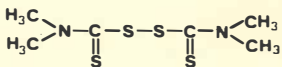
| lfd. Nr. | BBA Nr. | CIPAC Nr. | Name         | Chemische Bezeichnung  | Strukturformel   | Summenformel      | Mol.-gew.                  | Wir-<br>kungs-<br>bereich | weitere<br>Namen |
|----------|---------|-----------|--------------|--|--|-------------------|----------------------------|---------------------------|------------------|
| 241      | 392     | 268       | Resmethrin   | 5-Benzyl-3-furyl-<br>methyl-(±)-cis,<br>trans-chrysanthemat  |  | $C_{22}H_{26}O_3$ | 338.0                      | I                         |                  |
| 242      | 193     | 38        | Rotenon      | 1,2,12,12a-Tetra-<br>hydro-2-isopro-<br>penyl-8,9-di-<br>methoxy-(1)-benzo-<br>pyrano-(3,4-b)-<br>furo-(2,3-h)-(1)-<br>benzopyran-6-on |  | $C_{23}H_{22}O_6$ | 394.4                      | I<br>A                    |                  |
| 243      | 313     |           | S 421        | Di-(2,3,3,3-tetrachlor-<br>propyl)ether  | $Cl_3C-CH(Cl)-CH_2-O-CH_2-CH(Cl)-CCl_3$  | $C_6H_6Cl_8O$     | 377.7                      | S                         |                  |
| 244      | 184     | 18        |              | Schwefel   |  |                   | 256.5<br>(S <sub>8</sub> ) | F<br>(A)                  |                  |
| 245      | 252     |           | Sebuthylazin | 4-Ethylamino-2-sec.<br>butylamino-6-chlor-<br>1,3,5-triazin  |  | $C_9H_{16}ClN_5$  | 229.7                      | H                         |                  |
| 246      | 101     | 22        | Simazin      | 6-Chlor-2,4-bis-ethyl-<br>amino-1,3,5-triazin  |  | $C_7H_{12}ClN_5$  |                            |                           |                  |

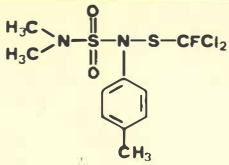
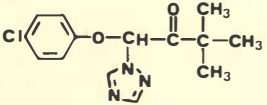
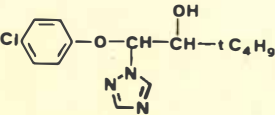
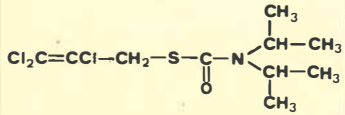
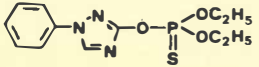


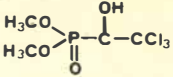
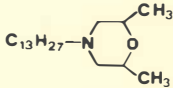
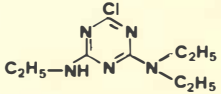
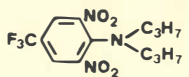
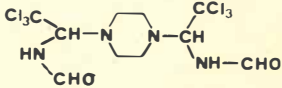
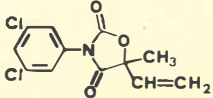
| lfd. Nr. | BBA Nr. | CIPAC Nr. | Name            | Chemische Bezeichnung                               | Strukturformel   | Summenformel          | Mol.-gew. | Wir-<br>kungs-<br>bereich | weitere<br>Namen |
|----------|---------|-----------|-----------------|---|--|-----------------------|-----------|---------------------------|------------------|
| 247      | 329     | 211       | Sulfachinoxalin | N-(Chinoxalin-2-yl)-sulfanilamid                    |  | $C_{14}H_{12}N_4O_2S$ | 300.3     | R                         |                  |
| 248      | 104     | 198       | Sulfotep        | O,O,O,O-Tetraethylthiodiphosphat                    |   | $C_8H_{20}O_5P_2S_2$  | 322.3     | A<br>I                    |                  |
| 249      | 105     | 6         | 2.4.5-T         | 2,4,5-Trichlorphenoxyessigsäure                     |   | $C_6H_5Cl_3O_3$       | 255.5     | H                         |                  |
| 250      | 107     | 67        | TCA             | Trichloressigsäure                                  | $Cl_3C-COOH$   | $C_2HCl_3O_2$         | 163.4     | H                         |                  |
| 251      | 417     | 272       | Terbacil        | 3-tert.-Butyl-5-chlor-6-methyluracil                |   | $C_9H_{13}ClN_2O_2$   | 216.7     | H                         |                  |
| 252      | 459     |           | Terbufos        | S-tert.-Butylthio-methyl-O,O-diethyl-dithiophosphat |  | $C_9H_{21}O_2PS_3$    | 288.4     | I                         |                  |

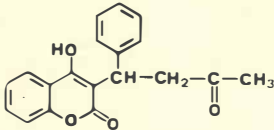
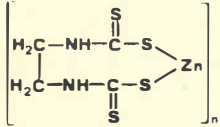
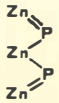
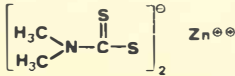
| Ifd. Nr. | BBA Nr. | CIPAC Nr. | Name              | Chemische Bezeichnung  | Strukturformel   | Summenformel  | Mol.-gew. | Wirkungsbereich | weitere Namen |
|----------|---------|-----------|-------------------|--|--|---|-----------|-----------------|---------------|
| 253      | 293     |           | Terbumeton        | 4-Ethylamino-2-tert.-butylamino-6-methoxy-1,3,5-triazin                          |  | C <sub>10</sub> H <sub>19</sub> N <sub>5</sub> O                | 225.3     | H               |               |
| 254      | 316     | 234       | Terbutylazin      | 4-Ethylamino-2-tert.butylamino-6-chlor-1,3,5-triazin                             |  | C <sub>9</sub> H <sub>16</sub> ClN <sub>5</sub>                 | 229.7     | H               |               |
| 255      | 246     | 212       | Terbutryn         | 4-Ethylamino-2-tert.-butylamino-6-methyl-thio-1,3,5-triazin                      |  | C <sub>10</sub> H <sub>19</sub> N <sub>5</sub> S                | 241.4     | H               |               |
| 256      | 317     | 265       | Tetrachlorvinphos | O-[2-Chlor-1-(2,4,5-trichlorphenyl)-vinyl]-O,O-dimethylphosphat (trans-Isomeres) |   | C <sub>10</sub> H <sub>9</sub> Cl <sub>4</sub> O <sub>4</sub> P | 366.0     | I               |               |
| 257      | 109     | 113       | Tetradifon        | 4-Chlorphenyl-2,4,5-trichlorphenylsulfon   |   | C <sub>12</sub> H <sub>6</sub> Cl <sub>4</sub> O <sub>2</sub> S | 356.1     | A               |               |

| lfd. Nr. | BBA Nr. | CIPAC Nr. | Name                      | Chemische Bezeichnung  | Strukturformel   | Summenformel       | Mol.-gew. | Wir-<br>kungs-<br>bereich | weitere<br>Namen |
|----------|---------|-----------|---------------------------|--|--|--------------------|-----------|---------------------------|------------------|
| 258      | 458     |           | Tetramethrin              | 3,4,5,6-Tetra-<br>hydrophthalimido-<br>methylchrysanthemat                     |  | $C_{19}H_{25}NO_4$ | 331.4     | I                         |                  |
| 259      | 110     | 114       | Tetrasul                  | 4'-Chlorphenyl-2,4,5-<br>trichlorphenylsulfid                                  |   | $C_{12}H_6Cl_4S$   | 324.1     | A                         |                  |
| 260      | 129     | 181       | Thallium-<br>verbindungen |  |  |                    |           | R                         |                  |
| 261      | 256     | 323       | Thiabendazol              | 2-(4-Thiazolyl)-<br>benzimidazol   |   | $C_{10}H_7N_3S$    | 201.3     | F                         |                  |
| 262      | 384     |           | Thiazfluron               | 1,3-Dimethyl-1-<br>(5-trifluor-methyl-<br>1,3,4-thiadiazol-<br>2-yl)-harnstoff |  | $C_6H_7F_3N_4O_3S$ | 240.2     | H                         |                  |
| 497      |         |           | Thiocyclam                | N,N-Dimethyl-1,2,3-<br>trithian-5yl-amin                                       |   | $C_5H_{10}NS_3$    | 180.3     | I                         |                  |

| Ifd. Nr. | BBA Nr. | CIPAC Nr. | Name              | Chemische Bezeichnung   | Strukturformel   | Summenformel            | Mol.-gew. | Wir-<br>kungs-<br>bereich | weitere<br>Namen |
|----------|---------|-----------|-------------------|---|--|-------------------------|-----------|---------------------------|------------------|
| 263      | 493     | 351       | Thiofanox         | 2,2-Dimethyl-1-methylthiomethyl-propylidenamino-methylcarbamat  |  | $C_9H_{18}N_2O_2S$      | 218.3     | I                         |                  |
| 264      | 319     | 115       | Thiometon         | S-[2-(Ethylthio)-ethyl]-O,O-dimethyl-dithiophosphat   |  | $C_6H_{15}O_2PS_3$      | 246.4     | I                         |                  |
| 265      | 204     | 182       | Thionazin         | O,O-Diethyl-O-(pyrazin-2-yl)thio-phosphat   |   | $C_8H_{13}N_2O_3PS$     | 248.2     | N                         |                  |
| 266      | 370     | 262       | Thiophanat-methyl | 1,2-Di-(3-methoxy-carbonyl-2-thio-ureido)benzol<br>oder<br>Dimethyl-4,4-O-phenylen-bis-(3-thioallophanat) |  | $C_{12}H_{14}N_4O_4S_2$ | 342.4     | F                         |                  |
| 267      | 119     | 24        | Thiram            | Tetramethylthiuram-disulfid   |  | $C_6H_{12}N_2S_4$       | 240.4     | F<br>P                    | TMTD             |

| lfd. Nr. | BBA Nr. | CIPAC Nr. | Name        | Chemische Bezeichnung   | Strukturformel   | Summenformel   | Mol.-gew. | Wirkungsbereich | weitere Namen |
|----------|---------|-----------|-------------|---|--|--|-----------|-----------------|---------------|
| 268      | 371     | 275       | Tolyfluamid | N-Dichlorfluor-methylthio-N',N'-dimethyl-N-p-tolyl-sulfamid       |   | C <sub>10</sub> H <sub>13</sub> Cl <sub>2</sub> FN <sub>2</sub><br>O <sub>2</sub> S <sub>2</sub> | 347.3     | F               |               |
| 269      | 425     | 352       | Triadimefon | 1-(4-Chlorphenoxy)-3,3-dimethyl-1-(1,2,4-triazol-1-yl)butan-2-on  |  | C <sub>14</sub> H <sub>16</sub> ClN <sub>3</sub> O <sub>2</sub>                                  | 293.8     | F               |               |
| 269a     | 605     |           | Triadimenol | 1-(4-Chlorphenoxy)-3,3-dimethyl-1-(1,2,4-triazol-1-yl)-butan-2-ol |  | C <sub>14</sub> H <sub>18</sub> ClN <sub>3</sub> O <sub>2</sub>                                  | 285.8     | F               |               |
| 270      | 135     | 97        | Tri-allat   | S-2,3,3-Trichlor-allyl-N,N-diisopropylthiocarbamat                |  | C <sub>10</sub> H <sub>16</sub> Cl <sub>3</sub> NOS  | 304.7     | H               |               |
| 271      | 401     |           | Triazophos  | O,O-Diethyl-O-(1-phenyl-1,2,4-triazol-3-yl)-thio-phosphat         |  | C <sub>12</sub> H <sub>16</sub> N <sub>3</sub> O <sub>3</sub> PS                                 | 313.3     | A<br>I<br>N     |               |

| Ifd. Nr. | BBA Nr. | CIPAC Nr. | Name        | Chemische Bezeichnung  | Strukturformel   | Summenformel             | Mol.-gew. | Wir-<br>kungs-<br>bereich | weitere<br>Namen |
|----------|---------|-----------|-------------|--|--|--------------------------|-----------|---------------------------|------------------|
| 272      | 112     | 68        | Trichlorfon | O,O-Dimethyl-2,2,2-trichlor-1-hydroxyethylphosphonat             |   | $C_4H_8Cl_3O_4P$         | 257.4     | I                         |                  |
| 273      | 320     | 324       | Tridemorph  | 2,6-Dimethyl-4-tridecyl-morpholin                                |   | $C_{19}H_{39}NO$         | 297.5     | F                         |                  |
| 274      | 386     |           | Trietazin   | 4-Ethylamino-6-chlor-2-diethyl-amino-1,3,5-triazin               |   | $C_9H_{15}ClN_5$         | 228.7     | H                         |                  |
| 275      | 321     | 183       | Trifluralin | 2,6-Dinitro-N,N-dipropyl-4-trifluor-methylanilin                 |   | $C_{13}H_{16}F_3N_3O_4$  | 335.3     | H                         |                  |
| 276      | 338     |           | Triforin    | 1,4-Di-(2,2,2-tri-chlor-1-formamido-ethyl)-piperazin             |  | $C_{10}H_{14}Cl_6N_4O_2$ | 434.9     | A<br>F                    |                  |
| 277      | 412     | 280       | Vinclozolin | 3-(3,5-Dichlor-phenyl)-5-methyl-5-vinyl-1,3-oxazoli-din-2,4-dion |   | $C_{12}H_9Cl_2NO_3$      | 286.1     | F                         |                  |

| lfd. Nr. | BBA Nr. | CIPAC Nr. | Name     | Chemische Bezeichnung                               | Strukturformel   | Summenformel        | Mol.-gew. | Wirkungsbereich | weitere Namen |
|----------|---------|-----------|----------|---|--|---------------------|-----------|-----------------|---------------|
| 278      | 374     |           | Vondozeb | Maneb-Zineb-Mischfällung mit 14,5 % Mn und 2,1 % Zn |  |                     |           | F               | Triziman D    |
| 279      | 114     | 70        | Warfarin | 4-Hydroxy-3-(3-oxo-1-phenylbutyl) coumarin          |  | $C_{19}H_{16}O_4$   | 308.2     | R               |               |
| 280      | 116     | 25        | Zineb    | Zink-ethylen-1,2-bis-dithiocarbamat                 |   | $C_4H_6N_2S_4Zn$    | 275.7     | F               |               |
| 281      | 3       | 69        |          | Zinkphosphid  |   | $Zn_3P_2$           | 258.1     | R               |               |
| 282      | 118     | 31        | Ziram    | Zink-dimethyldithiocarbamat                         |   | $C_6H_{12}N_2S_4Zn$ | 305.8     | F               |               |

