



AGROLAB LUFA Dr.-Hell-Str. 6, 24107 Kiel

COPELAND CROP SRL
Str 601, Nr. 54
317235 Pecica
RUMÄNIEN

Date 23.11.2021

Customer no. 10085356

REPORT 2992615 - 371339

Order 2992615
 Sample no. 371339
 Sample acceptance 17.11.2021
 Date of sampling 15.11.2021
 Sample taker Client
 Customer sample description Organic Corn
 Packaging 1x Quantum protect
 LOT-No./Batch LOT ORG 21

Unit

Result Declaration

Substance Method

Pesticides Multi-Residue-Methods (complete list see appendix)**In the range of performed analysis no pesticides were detected above limit of quantification.****Further sample data**

Amount of sample received	g	892			OM	no information
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Mycotoxins

Aflatoxine B1	µg/kg	<0,3			OM	QMP_504_KI_52_151 : 2020-11 (LC-MSMS)
Aflatoxine B2	µg/kg	<1,0 ^{mf)}			OM	QMP_504_KI_52_151 : 2020-11 (LC-MSMS)
Aflatoxine G1	µg/kg	<0,5			OM	QMP_504_KI_52_151 : 2020-11 (LC-MSMS)
Aflatoxine G2	µg/kg	<0,5			OM	QMP_504_KI_52_151 : 2020-11 (LC-MSMS)
Deoxynivalenol (DON)	µg/kg	404			OM	QMP_504_KI_52_151 : 2020-11 (LC-MSMS)
Zearalenone (ZEA)	µg/kg	<15 ^{mf)}			OM	QMP_504_KI_52_151 : 2020-11 (LC-MSMS)
Ochratoxine A	µg/kg	<0,500			OM	QMP_504_KI_52_151 : 2020-11 (LC-MSMS)
HT-2 Toxin	µg/kg	31,6			OM	QMP_504_KI_52_151 : 2020-11 (LC-MSMS)
T-2 Toxin	µg/kg	32,5			OM	QMP_504_KI_52_151 : 2020-11 (LC-MSMS)

GMO-Screening

GMO Screening CTP2-CP4-EPSPS genetic construct		not detected			OM	ASU L 00.00-154 : 2014-08 (mod.)
GMO Screening FMV-Promotor		not detected			OM	ASU L 00.00-148 : 2014-02 (mod.)
GMO Screening nos-Terminator		not detected			OM	ASU L 00.00-122 : 2008-06 (mod.)
GMO Screening pat-Gen		not detected			OM	ASU L 00.00-154 : 2014-08 (mod.)
GMO Screening 35S CaMV Promotor		not detected			OM	ASU L 00.00-122 : 2008-06 (mod.)

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x) The sum calculation is done without taking into account single values below limit of detection or limit of quantification.
wf) In the present sample the recovery of one or more internal standards is < 50% but > 10%. Consequently a higher measurement uncertainty is expected.

Explanation: The symbol "<" or n.d. in the result column means, the substance concerned is not quantifiable at the limit of quantification shown opposite.

The sign "<"...."(LOD)" or n.d. in column result means, the substance concerned cannot be detected within the limit of detection.

Parameter-specific analytical measurement uncertainties and information regarding the method of calculation will be provided upon request if the reported results are above the parameter-specific limit of quantification.

Explanation: OM = on original matter; DM = on dry matter base

The sampling date is a customer information.

Remark to amount of sample received: Total amount including packaging

Remarks

Start of testing: 17.11.2021

End of testing: 23.11.2021

The results are related only to the samples tested. In cases where the laboratory has not been responsible for sampling, the reported results apply to the samples as received. Duplication of this document or of parts of it requires the authorization from laboratory. In accordance our agreement in writing in the order confirmation, the results in this test report are in a simplified form in the context of DIN EN ISO/IEC 17025:2018, paragraph 7.8.1.3.



AGROLAB LUFA Frau Erika Paschke, Tel. 0431/1228-318
officially approved foodchemist
customer relation management food



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List of all analyzed pesticides (limit of quantification [mg/kg])

Method: calculated, Unit: mg/kg				
Parameter	Limit of quantification	Parameter	Limit of quantification	Limit of quantification
Sum acibenzolar-S-methyl and acibenzolar acid (without hydrolysis)		Sum aldicarb/-sulfon/-sulfoxid	Sum aldrin, dieldrin	
Sum amitraz		Sum bentazone	Sum bifenazate	
Sum captan and Tetrahydrophthalimide (THPI)		Sum carbofuran, 3-hydroxycarbofuran	Sum carboxin	
Sum chloridazon		Sum chlorpyrifos-methyl	Sum clethodim	
Sum cycloxydim		Sum DDT-isomers	Sum disulfoton	
Sum endosulfan-alpha, -beta, -sulfat		Sum ethofumesate	Sum fenamiphos, -sulphoxide, -sulphone	
Sum fenchlorphos		Sum fenthion	Sum fipronil, -sulfone (MB 46136)	
Sum flonicamid		Sum flufenacet	Sum heptachlor, heptachlorepoxyde	
Sum Isoxaflutole		Sum MCPA, MCPB (without hydrolysis)	Sum metazachlor	
Sum methiocarb, -sulfone, -sulfoxide		Sum milbemectin	Sum of cis- and trans-chlordane (F) (R)	
Sum of Folpet and Phthalimid		Sum of malathion and malaoxon	Sum oxydemeton-methyl, demeton-S-methyl-sulfon	
Sum Parathion-methyl		Sum phorate	Sum phosmet and phosmet-oxon	
Sum prochloraz		Sum propachlor	Sum propoxycarbazone	
Sum pyridate (without hydrolysis)		Sum quintozone and pentachloro-aniline	Sum spirotetramat	
Sum tepraloxymid		Sum tolylfluand	Sum triflumizole and FM 6-1	
1-naphthylacetamide and 1-naphthylacetic acid				

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Method: EN 15662 : 2018 (mod.), Unit: mg/kg					
Parameter	Limit of quantification	Parameter	Limit of quantification	Parameter	Limit of quantification
Accephate	0,01	Acetamidiprid	0,01	Acibenzolaracid (free acid)	0,01
Acibenzolar-S-methyl (before hydrolysis)	0,01	Acronifen	0,01	Acrinathrin and its enantiomer	0,01
Alachlor	0,01	Aldicarb	0,01	Aldicarb-sulfon	0,01
Aldicarb-sulfoxide	0,01	Aldrin	0,005	Ametoctradin	0,01
Ametryn	0,01	Aminocarb	0,01	Amitraz	0,01
Anthraquinone	0,01	Atrazine	0,01	Azaconazole	0,01
Azadirachtin	0,01	Azinphos-ethyl	0,01	Azinphos-methyl	0,01
Azoxystrobin	0,01	Benalaxyl	0,01	Bendiocarb	0,01
Benfluralin	0,01	Bensulfuron-methyl	0,01	Bentazone	0,01
Benthiavalicarb-isopropyl	0,01	Benzovindiflupyr	0,01	Bifenazate	0,01
Bifenazate-diazene	0,01	Bifenox	0,01	Bifenthrin	0,01
Biphenyl	0,01	Bitertanol	0,01	Bixafen	0,01
Boscalid	0,01	Bromacil	0,01	Bromocyclen	0,01
Bromophos-ethyl	0,01	Bromophos-methyl	0,01	Bromopropylate	0,01
Bromoxynil	0,01	Bromuconazole	0,01	Bupirimate	0,01
Buprofezin	0,01	Butafenacil	0,01	Butocarboxim	0,01
Butocarboxim-sulfoxide	0,01	Butoxycarboxim	0,01	Cadusafos	0,01
Captan	0,01	Carbaryl	0,01	Carbofuran	0,01
Carbophenothion	0,01	Carbophenothion-methyl	0,01	Carbosulfan	0,01
Carboxin	0,01	Chlorantraniliprol	0,01	Chlorbenside	0,01
Chlorbufam	0,01	Chlordane alpha	0,005	Chlordane gamma	0,005
Chlordane oxy	0,005	Chlorfenapyr	0,01	Chlorfenprop-methyl	0,01
Chlorfenson	0,01	Chlorfluazuron	0,01	Chlorflurenol	0,01
Chlorflurenol-methyl	0,01	Chloridazon	0,01	Chlorimuron-ethyl	0,01
Chlormephos	0,01	Chlorobenzilate	0,01	Chloroneb	0,01
Chlorotoluron	0,01	Chlorphenvinphos	0,01	Chlorpropham	0,01
Chlorpropylate	0,01	Chlorpyrifos	0,01	Chlorpyrifos-methyl	0,01
Chlorpyrifos-methyl-desmethyl	0,01	Chlorthal-dimethyl	0,01	Chlorthalonil	0,01
Chlorthion	0,01	Chlorthiophos	0,01	Chlozolate	0,01
Cinosulfuron	0,01	Clethodim	0,01	Climbazole	0,01
Clodinafop	0,01	Clodinafop-propargyl	0,01	Clofentezin	0,01
Clomazone	0,01	Cloquintocet-mexyl	0,01	Clothianidin	0,01
Coumaphos	0,01	Crimidine	0,01	Cyanazin	0,01
Cyanofenphos	0,01	Cyanophos	0,01	Cyantraniliprol	0,01
Cyazofamid	0,01	Cyclanilid	0,01	Cycloate	0,01
Cycloxydim	0,01	Cyflufenamid	0,01	Cyflumetofen	0,01
Cyfluthrin	0,01	Cyhalofop-butyl	0,01	Cyhalothrine	0,01
Cymoxanil	0,01	Cypermethrin	0,01	Cyproconazole	0,01
Cyprodinil	0,01	Deltamethrin	0,01	Demeton-S-methyl	0,01
Demeton-S-methyl-sulfone	0,01	Desethylatrazine	0,01	Desmedipham	0,01
Desmetryn	0,01	Diazinon	0,01	Dichlobenil	0,01
Dichlofenthione	0,01	Dichlofluanid	0,01	Dichlorprop (free acid)	0,01
Dichlorvos	0,01	Diclobutrazole	0,01	Diclofop	0,01
Dicloran	0,01	Dicofol	0,01	Dicrotophos	0,01
Dieldrin	0,005	Diethofencarb	0,01	Diethyltoluamide (DEET)	0,01
Difenacoum	0,01	Difenoconazole	0,01	Diflubenzuron	0,01
Diflufenican	0,01	Dimethenamide	0,01	Dimethoate	0,01
Dimethomorph	0,01	Dimethylaminosulfotoluidide (DMST)	0,01	Dimoxystrobin	0,01
Diniconazole	0,01	Dinocap	0,01	Dinotefuran	0,01
Dinoterb (before hydrolysis)	0,01	Diphenamid	0,01	Diphenylamine	0,01
Dipropetryn	0,01	Disulfoton	0,01	Disulfoton-sulfone	0,01
Disulfoton-sulfoxide	0,01	Ditalimfos	0,01	Diuron	0,01
DMSA	0,01	Dodemorph	0,01	Dodin	0,01
Emamectin	0,01	Endosulfan alpha	0,005	Endosulfan beta	0,005
Endosulfansulfat	0,005	Endrin	0,005	Endrin Ketone	0,01
EPN	0,01	Epoxiconazole	0,01	EPTC	0,01
Etaconazole	0,01	Ethalfuralin	0,01	Ethiofencarb	0,01
Ethiofencarb-sulfon	0,01	Ethiofencarb-sulfoxide	0,01	Ethion	0,01
Ethiprole	0,01	Ethirimol	0,01	Ethofumesate	0,01
Ethofumesate-2-keto	0,05	Ethoprophos	0,01	Ethoxyquin	0,01
Etofenprox	0,01	Etoxazole	0,01	Etridiazole	0,01
Etrifos	0,01	Famoxadone	0,01	Famphur	0,01
Fenamidone	0,01	Fenamiphos	0,01	Fenamiphos-sulfoxide	0,01
Fenamiphos-sulphone	0,01	Fenarimole	0,01	Fenazaquine	0,01
Fenbuconazole	0,01	Fenbutatin oxide	0,01	Fenchlorphos	0,01
Fenchlorphos-oxon	0,01	Fenfluthrin	0,01	Fenhexamid	0,01

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Parameter	Limit of quantification	Parameter	Limit of quantification	Parameter	Limit of quantification
Fenitrothion	0,01	Fenobucarb	0,01	Fenoxaprop	0,01
Fenoxycarb	0,01	Fenpiclonil	0,01	Fenpropathrine	0,01
Fenpropidin	0,01	Fenpropimorph	0,01	Fenpyrazamin	0,01
Fenpyroximate	0,01	Fenson	0,01	Fensulfothion	0,01
Fensulfothion-oxon	0,01	Fensulfothion-oxon-sulfon	0,01	Fensulfothion-sulfon	0,01
Fenthion	0,01	Fenthion-oxone	0,01	Fenthion-oxon-sulfon	0,01
Fenthionoxonsulfoxide	0,01	Fenthion-sulfon	0,01	Fenthion-sulfoxide	0,01
Fentin	0,01	Fenuron	0,01	Fenvalerate	0,01
Fipronil	0,002	Fipronil-sulfon	0,002	Flonicamid	0,01
Fluazifop (free acid)	0,01	Fluazifop-butyle	0,01	Fluazinam	0,01
Flubendiamid	0,01	Fluchloralin	0,01	Flucythrinate	0,01
Fludioxonil	0,01	Flufenacet	0,01	Flufenacet ESA (ethansulfonic acid)	0,01
Flufenacet OA (Oxalamic Acid)	0,01	Flufenacet-alcohol	0,01	Flufenoxuron	0,01
Flufenzin	0,01	Flumetralin	0,01	Flumioxazin	0,01
Fluometuron	0,01	Flupicolide	0,01	Fluopyram	0,01
Fluquinconazole	0,01	Flurochloridone	0,01	Fluroxypyr (free acid)	0,01
Flurprimidol	0,01	Flusilazole	0,01	Fluthiacet-methyl	0,01
Flutolanil	0,01	Flutriafof	0,01	Fluxapyroxad	0,01
FM 6-1	0,01	Folpet	0,01	Fonofos	0,01
Forchlorfenuron	0,01	Formetanate(hydrochloride)	0,01	Formothion	0,01
Fosthiazat	0,01	Fuberidazole	0,01	Furalaxyl	0,01
Furathiocarb	0,01	Genite	0,01	Halfenprox	0,01
Halofenozid	0,01	Haloxypop (free acid)	0,01	Haloxypop methyl	0,01
Haloxypop-ethoxy-ethyl	0,01	HCH-alpha	0,005	HCH-beta	0,005
HCH-delta	0,005	HCH-epsilon	0,005	HCH-gamma (Lindane)	0,005
Heptachlor	0,005	Heptachlorepoxyde-cis	0,005	Heptachlorepoxyde-trans	0,005
Heptenophos	0,01	Hexachlorobenzene	0,005	Hexaconazole	0,01
Hexaflumuron	0,01	Hexazinone	0,01	Hexithiazox	0,01
Icaridin (Picaridin)	0,01	Imazalil	0,01	Imazamox	0,01
Imazapic	0,01	Imazaquine	0,01	Imazethapyr	0,01
Imibenconazole	0,01	Imidacloprid	0,01	Indoxacarb	0,01
Iodofenphos	0,01	Iodosulfuron-methyl-sodium	0,01	Ioxynil	0,01
Iprobenfos	0,01	Iprodion	0,01	Iprovalicarb	0,01
Isazofos	0,01	Isocarbophos	0,01	Isodrin	0,01
Isofenphos	0,01	Isofenphos-methyl	0,01	Isofetamid	0,01
Isoprocab	0,01	Isoprothiolane	0,01	Isoproturon	0,01
Isoprazam	0,01	isoxaben	0,01	Isxadifen-ethyl	0,01
Isoxaflutole	0,01	Isoxathion	0,01	Kresoxim-methyl	0,01
lambda-cyhalothrin	0,01	Landrin (3,4,5-Trimethacarb)	0,01	Lenacil	0,01
Leptophos	0,01	Linuron	0,01	Malafoxon	0,01
Malathion	0,01	Mandestrobin	0,01	Mandipropamid	0,01
MCPA (free acid)	0,01	MCPB (free acid)	0,01	Mecarbame	0,01
Mecoprop	0,01	Mefenpyr-diethyl	0,01	Mepanipyrim	0,01
Mepronil	0,01	Meptyldinocap	0,01	Metaflumizone	0,01
Metalaxyl (Sum of Metalaxyl and Metalaxyl-M)	0,01	Metamitron	0,01	Metazachlor	0,01
Metconazole	0,01	Methabenzthiazuron	0,01	Methacrifos	0,01
Methamidophos	0,01	Methidathion	0,01	Methiocarb	0,01
Methiocarb-sulfon	0,01	Methiocarb-sulfoxid	0,01	Methomyl	0,01
Methoprotyrine	0,01	Methoxychlor	0,005	Methoxyfenozide	0,01
Metobromuron	0,01	Metolachlor	0,01	Metolcarb	0,01
Metosulam	0,01	Metoxuron	0,01	Metrafenone	0,01
Metribuzin	0,01	Metsulfurone-methyl	0,01	Mevinphos	0,01
Milbemectin A3	0,01	Milbemectin A4	0,01	Mirex	0,005
Molinate	0,01	Monocrotophos	0,01	Monolinuron	0,01
Monuron	0,01	Myclobutanil	0,01	Napropamide	0,01
Neburon	0,01	Nicosulfuron	0,01	Nitenpyram	0,01
Nitratin	0,01	Nitrapyrin	0,01	Nitrofen	0,005
Nitrothal-isopropyl	0,01	Norflurazone	0,01	Novaluron	0,01
Nuarimol	0,01	N-2,4-Dimethylphenyl-N-methylformamidine	0,01	Octachlordipropylether (S421)	0,01
Ofurace	0,01	Omethoate	0,01	o,p-DDD	0,005
o,p-DDE	0,005	o,p-DDT	0,005	Oxadiazon	0,01
Oxadixyle	0,01	Oxamyl	0,01	Oxamyl-oxim	0,01
Oxydemeton-methyl	0,01	Oxyfluorfen	0,01	Paclobutrazol	0,01
Paraoxon-ethyl	0,01	Paraoxon-methyl	0,02	Parathion-ethyl	0,01
Parathion-methyl	0,01	Pebulate	0,01	Penconazol	0,01

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Parameter	Limit of quantification	Parameter	Limit of quantification	Parameter	Limit of quantification
Pencycuron	0,01	Pendimethalin	0,01	Pentachloro-aniline	0,01
Pentachloroanisol	0,01	Pentachlorobenzene	0,01	Penthiopyrad	0,01
Permethrin	0,01	Perthane	0,01	Pethoxamid	0,01
Phenkapton	0,01	Phenmedipham	0,01	Phenthoate	0,01
Phorate	0,01	Phorat-oxon	0,01	Phorat-oxon-sulfon	0,01
Phorat-oxon-sulfoxid	0,01	Phorat-sulfon	0,01	Phorat-sulfoxid	0,01
Phosalone	0,01	Phosmet	0,01	Phosmet-oxon	0,01
Phosphamidon	0,01	Phthalimide	0,02	Picolinafen	0,01
Picoxystrobin	0,01	Piperonylbutoxide	0,01	Pirimicarb	0,01
Pirimicarb-desmethyl	0,01	Pirimiphos-ethyl	0,01	Pirimiphos-methyl	0,01
p,p-DDD	0,005	p,p-DDE	0,005	p,p-DDT	0,005
Prochloraz	0,01	Prochloraz desimidazole-amino (BTS 44595)	0,01	Prochloraz desimidazole-formylamino (BTS 44596)	0,01
Procymidone	0,01	Profenofos	0,01	Profluralin	0,01
Profoxydim	0,01	Promecarb	0,01	Prometryn	0,01
Propachlor	0,01	Propachlor OA (Oxalamic Acid)	0,01	Propamocarb	0,01
Propanil	0,01	Propaquizafop	0,01	Propargite	0,01
Propazine	0,01	Propetamphos	0,01	Propham	0,01
Propiconazole	0,01	Propoxur	0,01	Propoxy-carbazone	0,01
Propyzamide	0,01	Proquinazide	0,01	Prosulfocarb	0,01
Prothioconazole (Prothioconazole-desthio)	0,01	Prothiophos	0,01	Pymetrozine	0,01
Pyraclostrobin	0,01	Pyraflufen-ethyl	0,01	Pyrazophos	0,01
Pyrethrins	0,01	Pyridaben	0,01	Pyridalyl	0,01
Pyridaphenthion	0,01	Pyridate (without hydrolysis)	0,01	Pyrifenox	0,01
Pyrimethanil	0,01	Pyrimidifen	0,01	Pyriproxyfen	0,01
Pyroxsulam	0,01	Quinalphos	0,01	Quinoxyfen	0,01
Quintozene	0,005	Quizalofop (free acid)	0,01	Quizalofop-ethyl	0,01
Resmethrine	0,01	Rotenone	0,01	Sedaxane	0,01
Sethoxydim	0,01	Silafluofen	0,01	Silthiofam	0,01
Simazin	0,01	Spinetoram	0,01	Spinosad	0,01
Spiromesifen	0,01	Spirotetramat	0,01	Spirotetramat-enol	0,01
Spirotetramat-enol-glucoside	0,01	Spirotetramat-ketohydroxy	0,01	Spirotetramat-monohydroxy	0,01
Spiroxamine	0,01	Sulfentazone	0,01	Sulfotep	0,01
Sulfoxaflor	0,01	Sulprofos	0,01	Sum carbendazim/benomyl	0,01
tau-Fluvalinate	0,01	Tebuconazole	0,01	Tebufenozide	0,01
Tebufenpyrad	0,01	Tecnazene	0,005	Teflubenzuron	0,01
Tefluthrine	0,01	Tembotrion	0,01	Tepraloxymid	0,01
Terbacil	0,01	Terbufos	0,01	Terbufos-sulfon	0,01
Terbufos-sulfoxide	0,01	Terbumeton	0,01	Terbutryne	0,01
Terbutylazin-desethyl	0,01	Terbutylazine	0,01	Tetrachlorvinphos	0,01
Tetraconazole	0,01	Tetradifon	0,005	Tetrahydrophthalimide (THPI)	0,01
Tetrasul	0,01	TFNA	0,01	TFNG	0,01
Thiabendazole	0,01	Thiacloprid	0,01	Thiamethoxam	0,01
Thiobencarb	0,01	Thiodicarb	0,01	Thiofanox	0,01
Thiofanox-sulfoxide	0,01	Thiometon	0,01	Thiometon-sulfon	0,01
Thiometon-sulfoxid	0,01	Thiophanat-methyl	0,01	Tolclofos-methyl	0,01
Tolyfluanide	0,01	Tralkoxydim	0,01	Transfluthrine	0,01
Triadimefon	0,01	Triadimenol	0,01	Triallate	0,01
Triasulfuron	0,01	Triazamat	0,01	Triazophos	0,01
Trichlorfon	0,01	Trichloronate	0,01	Triclopyr	0,01
Tricyclazole	0,01	Tridemorph	0,01	Trifloxystrobin	0,01
Triflumizole	0,01	Triflumuron	0,01	Trifluralin	0,01
Triflusaluron-methyl	0,01	Triforine	0,01	Trinexapac	0,02
Trinexapac-ethyl	0,01	Triticonazole	0,01	Tritosulfuron	0,01
Uniconazole	0,01	Valifenalate	0,01	Vamidothion	0,01
Vinclozolin	0,01	Warfarin	0,01	Zoxamide	0,01
1-Naphthylacetic acid	0,05	1-Naphthylacetic amide	0,01	2-hydroxypropoxycarbazone	0,01
2-Naphthoxyacetic acid	0,01	2-Phenylphenol	0,01	2,4-D (free acid)	0,01
2,4-DB (free acid)	0,01	2,4-Dimethylphenylformamide	0,01	2,4,5-T (free acid)	0,01
3-Chloroaniline	0,01	3-Hydroxy-Carbofuran	0,01	4-Chlorophenoxyacetic acid (4-CPA)	0,01
4,4'-Dibromobenzophenone	0,01	6-hydroxy-Bentazone	0,01	8-hydroxy-Bentazone	0,01

wf) In the present sample the recovery of one or more internal standards is < 50% but > 10%. Consequently a higher measurement uncertainty is expected.

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Remark to 1-naphthylacetamide and 1-naphthylacetic acid: Sum of 1-naphthylacetamide and 1-naphthylacetic acid and its salts, expressed as 1-naphthylacetic acid. The sum parameter takes into account the active metabolites, which are detectable safely using the specified method. The actual content may be higher and can only be determined with a single method.

Remark to Benalaxyl: Benalaxyl including other mixtures of constituent isomers including Benalaxyl-M (sum of isomers).

Remark to Benthiavalicarb-isopropyl: Benthiavalicarb-isopropyl (KIF-230 R-L) and its enantiomer (KIF-230 S-D) and its diastereomers (KIF-230 S-L and KIF-230 R-D), expressed as benthiavalicarb-isopropyl (A).

Remark to Bifenthrin: Sum of isomers (F).

Remark to Bromoxynil: Bromoxynil and its salts, expressed as bromoxynil.

Remark to Bromuconazole: Sum of diastereoisomers (F).

Remark to Cyflufenamid: Sum of cyflufenamid (Z-isomer) and its E-isomer.

Remark to Cyfluthrin: Cyfluthrin including other mixtures of constituent isomers (sum of isomers) (F).

Remark to Cypermethrin: Cypermethrin including other mixtures of constituent isomers (sum of isomers) (F).

Remark to Diclofop: Sum diclofop-methyl and diclofop acid expressed as diclofop-methyl. By the multi-method only the free acid of the active ingredient is detected. If contents equal or higher than 0.008 mg/kg are detected, a quantitative analysis of the total acid is performed by hydrolysis.

Remark to Dicofof: Sum of p, p' and o, p' isomers (F).

Remark to Dimethenamid: Dimethenamid including other mixtures of constituent isomers including dimethenamid-P (sum of isomers).

Remark to Dimethomorph: Sum of isomers.

Remark to Diniconazole: Sum of isomers.

Remark to Dinocap: Sum of dinocap isomers and their corresponding phenols expressed as dinocap.

Remark to Fenpropidin: Sum of fenpropidin and its salts, expressed as fenpropidin (R) (A).

Remark to Fenpropimorph: Sum of isomers (F) (R).

Remark to Fentin: Fentin including its salts, expressed as triphenyltin cation (F).

Remark to Fenvaleate: Any ratio of constituent isomers (RR, SS, RS & SR) including esfenvaleate (F) (R).

Remark to Formetanate(hydrochloride): Sum of formetanate and its salts expressed as formetanate(hydrochloride).

Remark to HCH-alpha: Hexachlorocyclohexane (HCH), alpha-isomer (F).

Remark to HCH-beta: Hexachlorocyclohexane (HCH), beta-isomer (F).

Remark to HCH-gamma (Lindane): Lindane (Gamma-isomer of hexachlorocyclohexane (HCH)) (F).

Remark to Haloxypop-ethoxy-ethyl: By the multi-method only the free acid of the active ingredient is detected. If contents equal or higher than 0.008 mg/kg are detected, a quantitative analysis of the total acid is performed by hydrolysis.

Remark to Imazamox: Sum of imazamox and its salts, expressed as imazamox.

Remark to Indoxacarb: Sum of indoxacarb and its R enantiomer (F).

Remark to Iodosulfuron-methyl-sodium: Sum of iodosulfuron-methyl and its salts, expressed as iodosulfuron-methyl.

Remark to Ioxynil: Sum of ioxynil, its salts and its esters, expressed as ioxynil (F). By the multi-method only the free acid of the active ingredient is detected. If contents equal or higher than 0.008 mg/kg are detected, a quantitative analysis of the total acid is performed by hydrolysis.

Remark to Mecoprop: Sum of mecoprop-p and mecoprop expressed as mecoprop.

Remark to Meptyldinocap: Sum of 2,4 DNOPC and 2,4 DNOP expressed as meptyldinocap. By the multi-method only the free acid of the active ingredient is detected. If contents equal or higher than 0.008 mg/kg are detected, a quantitative analysis of the total acid is performed by hydrolysis.

Remark to Metaflumizone: Sum of E- and Z-isomers.

Remark to Metalaxyl (Sum of metalaxyl and metalaxyl-M): Metalaxyl including other mixtures of constituent isomers including metalaxyl-M (sum of isomers).

Remark to Metconazole: Sum of isomers (F).

Remark to Metolachlor: Metolachlor including other mixtures of constituent isomers including S-metolachlor (sum of isomers).

Remark to Mevinphos: Sum of E- and Z-isomers.

Remark to Paclobutrazol: Sum of the isomers.

Remark to Penconazole: Penconazole (Sum of isomers) (F).

Remark to Permethrin: Sum of isomers (F).

Remark to Propamocarb: Sum of propamocarb and its salts, expressed as propamocarb (R).

Remark to Propiconazole: Sum of the isomers (F).

Remark to Prothioconazole (Prothioconazole-desthio): Prothioconazole-desthio (sum of isomers) (F).

Remark to Resmethrin: Resmethrin including other mixtures of constituent isomers (sum of isomers) (F).

Remark to Spinosad: Spinosad, sum of spinosyn A and spinosyn D (F).

Remark to Spiroxamine: Sum of isomers (A) (R).

Remark to Sulfoxaflor: Sum of isomers.

Remark to Sum Amitraz: Amitraz including the metabolites containing the 2,4 -dimethylaniline moiety expressed as amitraz. The sum parameter takes into account the active metabolites, which are detectable safely using the specified method. The actual content may be higher and can only be determined with a single method.

Remark to Sum Carboxin: Carboxin (carboxin plus its metabolites carboxin sulfoxide and oxycarboxin (carboxin sulfone), expressed as carboxin). The sum parameter takes into account the active metabolites, which are detectable safely using the specified method. The actual content may be higher and can only be determined with a single method.

Remark to Sum DDT-isomers: Sum of p, p'-DDT, o, p'-DDT, p, p'-DDE and p, p'-TDE (DDD) expressed as DDT (F).

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Remark to Sum Flufenacet: Sum of all compounds containing the N fluorophenyl-N-isopropyl moiety expressed as flufenacet equivalent. The sum parameter takes into account the active metabolites, which are detectable safely using the specified method. The actual content may be higher and can only be determined with a single method.

Remark to Sum Isoxaflutole: Isoxaflutole (sum of isoxaflutole and its diketonitrile-metabolite, expressed as isoxaflutole)
Remark to Sum MCPA, MCPB: MCPA and MCPB (MCPA, MCPB including their salts, esters and conjugates expressed as MCPA) (R)
(F). The residue definition is not fully met as no hydrolysis has taken place in the multi-method.

Remark to Sum Milbemectin: Milbemectin (sum of milbemycin A4 and milbemycin A3 expressed as milbemectin).

Remark to Sum Pyridate: Sum of pyridate, its hydrolysis product CL 9673 (6-chloro-4-hydroxy-3-phenylpyridazin) and hydrolysable conjugates of CL 9673 expressed as pyridate).

The residue definition is not fully met as no hydrolysis has taken place in the multi-method.

Remark to Sum Spirotetramat: Summe aus Spirotetramat und Spirotetramat-enol, ausgedrückt als Spirotetramat (R)

Remark to Sum acibenzolar-S-methyl and acibenzolar: Sum of acibenzolar-S-methyl and acibenzolar acid (free and conjugated), expressed as acibenzolar-S-methyl. The residue definition is not fully met as no hydrolysis has taken place in the multi-method.

Remark to Sum aldicarb/-sulfon/-sulfoxid: Sum of aldicarb, its sulfoxide and its sulfone, expressed as aldicarb.

Remark to Sum aldrin, dieldrin: Aldrin and dieldrin combined expressed as dieldrin (F).

Remark to Sum bentazone: Sum of bentazone, its salts and 6-hydroxy (free and conjugated) and 8-hydroxy bentazone (free and conjugated), expressed as bentazone (R).

Remark to Sum bifenazate: Sum of bifenazate plus bifenazate-diazene expressed as bifenazate (F).

Remark to Sum captan and THPI: Sum of captan and THPI, expressed as captan (R) (A).

Remark to Sum carbendazim/benomyl: Sum of benomyl and carbendazim expressed as carbendazim (R).

Remark to Sum carbofuran, 3-hydroxycarbofuran: Sum of carbofuran (including any carbofuran generated from carbosulfan, benfuracarb or furathiocarb) and 3-OH carbofuran expressed as carbofuran (R).

Remark to Sum chloridazon: Chloridazon (R) (sum of chloridazon and chloridazon-desphenyl, expressed as chloridazon). The sum parameter takes into account the active metabolites, which are detectable safely using the specified method. The actual content may be higher and can only be determined with a single method.

Remark to Sum clethodim: Sum of sethoxydim and clethodim including degradation products calculated as sethoxydim. The sum parameter takes into account the active metabolites, which are detectable safely using the specified method. The actual content may be higher and can only be determined with a single method.

Remark to Sum cycloxydim: Cycloxydim including degradation and reaction products which can be determined as 3-(3-thianyl)glutaric acid S-dioxide (BH 517-TGSO₂) and/or 3-hydroxy-3-(3-thianyl)glutaric acid S-dioxide (BH 517-5-OH-TGSO₂) or methyl esters thereof, calculated in total as cycloxydim. The sum parameter takes into account the active metabolites, which are detectable safely using the specified method. The actual content may be higher and can only be determined with a single method.

Remark to Sum disulfoton: Sum of disulfoton, disulfoton sulfoxide and disulfoton sulfone expressed as disulfoton (F).

Remark to Sum endosulfan-alpha, -beta, -sulphate: Sum of alpha- and beta-isomers and endosulfan-sulphate expressed as endosulfan (F).

Remark to Sum ethofumesate: Sum of ethofumesate, 2-keto-ethofumesate, open-ring-2-keto-ethofumesate and its conjugate, expressed as ethofumesate. The sum parameter takes into account the active metabolites, which are detectable safely using the specified method. The actual content may be higher and can only be determined with a single method.

Remark to Sum fenamiphos, -sulfoxide, -sulfone: Sum of fenamiphos and its sulphoxide and sulphone expressed as fenamiphos.

Remark to Sum fenchlorphos: Sum of fenchlorphos and fenchlorphos oxon expressed as fenchlorphos.

Remark to Sum fipronil, -sulfone (MB 46136): Sum fipronil + sulfone metabolite (MB46136) expressed as fipronil (F).

Remark to Sum flonicamid: Sum of flonicamid, TFNA and TFNG expressed as flonicamid (R).

Remark to Sum folpet and phthalimide: Sum of folpet and phthalimide, expressed as folpet (R).

Remark to Sum heptachlor, heptachlorepoxy: Sum of heptachlor and heptachlor epoxide expressed as heptachlor (F).

Remark to Sum malathion and malaaxon: Sum of malathion and malaaxon expressed as malathion.

Remark to Sum metazachlor: Sum of metabolites 479M04, 479M08, 479M16, expressed as metazachlor (R). The sum parameter takes into account the active metabolites, which are detectable safely using the specified method. The actual content may be higher and can only be determined with a single method.

Remark to Sum methiocarb, -sulfone, -sulfoxide: Sum of methiocarb and methiocarb sulfoxide and sulfone, expressed as methiocarb.

Remark to Sum of cis- and trans-chlordane (F) (R): Chlordane (sum of cis- and trans-chlordane)

Remark to Sum oxydemeton-methyl, demeton-S-methyl-sulfon: Sum of oxydemeton-methyl and demeton-S-methylsulfone expressed as oxydemeton-methyl.

Remark to Sum parathion-methyl: Sum of Parathion-methyl and paraoxon-methyl expressed as Parathion-methyl.

Remark to Sum phorate: Sum of phorate, its oxygen analogue and their sulfones expressed as phorate.

Remark to Sum phosmet and phosmet-oxon: Phosmet and phosmet oxon expressed as phosmet (R).

Remark to Sum prochloraz: Sum of prochloraz and its metabolites containing the 2,4,6-Trichlorophenol moiety expressed as prochloraz.

Remark to Sum propachlor: Oxalinic derivate of propachlor, expressed as propachlor.

Remark to Sum propoxycarbazon: Propoxycarbazon, its salts and 2-hydroxypropoxycarbazon expressed as propoxycarbazon.

Remark to Sum quintozone and pentachloro-aniline: Sum of quintozone and pentachloro-aniline expressed as quintozone (F).

Remark to Sum tepaloxym: Sum of tepaloxym and its metabolites that can be hydrolysed either to the moiety 3-(tetrahydro-pyran-4-yl)-glutaric acid or to the moiety 3-hydroxy-(tetrahydro-pyran-4-yl)-glutaric acid, expressed as tepaloxym. The sum parameter takes into account the active metabolites, which are detectable safely using the specified method. The actual content may be higher and can only be determined with a single method.

Remark to Sum tolylfluorid: Sum of tolylfluorid and dimethylaminosulfotoluidide expressed as tolylfluorid (F) (R).

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Remark to Sum triflumizole and FM 6-1: Triflumizole and metabolite FM-6-1(N-(4-chloro-2-trifluoromethylphenyl)-n-propoxyacetamide), expressed as Triflumizole (F).

Remark to Tralkoxydim: Sum of the constituent isomers of tralkoxydim.

Remark to Trinexapac: Sum of trinexapac (acid) and its salts, expressed as trinexapac.

Remark to hydrolysis-relevant substances without carrying out the hydrolysis module: The validated limit of quantification is 0.01 mg/kg.

All data below this determination limit are to be interpreted as non-quantifiable traces. The actual content including the bound residues can only be determined via an additional hydrolysis step.

Remark to sum fenthion: Fenthion and its oxygen analogue, their sulfoxides and sulfone expressed as parent (F).

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