

KERAKOLL SpA Headquarter via Pedemontana, 25 41049 Sassuolo (MO) Italy

Test Report No. 59983-A010-AgBB-L

Test objective:

Proof of conformity with the AgBB scheme 2024

SKIL REMOVE (PTKK00302 | ID13800)

SKIL REMOVE (PTKK00302 | ID13800)

30/07/2025

Number of pages of report:

19

Testing / responsible laboratory:

eco-INSTITUT Germany GmbH, Köln

Note:

Test objective fulfilled:

The test results in the report refer exclusively to the test sample submitted by the manufacturer. The report is not permitted to be used in product and company advertising. The report may be published in full as technical documentation on the Internet with the written consent of eco-INSTITUT Germany GmbH. eco-INSTITUT Germany GmbH has recommended that the manufacturer repeats the test after 3 years at the latest. More information at www.eco-institut.de/en/advertising







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Sample View

Internal sample number (filled in by laboratory)

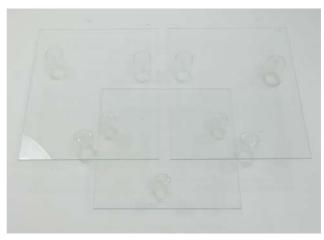
Photo of the test specimen: A010

Article designation according to order: Sample/batch number according to order: Type of sample:

Date of production: Sampling by: Date of sampling: Location of sampling:

Receipt of sample / Condition upon delivery:

59983-A010



SKIL REMOVE (PTKK00302 | ID13800)

PTKK00302

Waterbased restoring product for surfaces affected by mould, algae, fungi, lichen

15240605 06.2024

Marina Bastarolo

03/03/2025

no information

24/03/2025 / without objection



Statement of conformity with AgBB 2024

The sample with the internal sample number 59983-A010 has been tested on behalf of **KERAKOLL SpA Headquarter**. The article description according to the order is **SKIL REMOVE (PTKK00302 | ID13800)**.

This evaluation is based on the test criteria of the scheme "Health-related Evaluation of Emissions of Volatile Organic Compounds (VVOC, VOC and SVOC) from Building Products" of the Committee for Health-Related Evaluation of Building Products (AgBB 2024).

The results documented in the test report were evaluated as follows.¹

Test parameter	Result			Requirement		Requirement hold [yes/no]	
Emission analysis							
Measurement time: 3 days after test chamber loading							
Sum VOC (C6-C16) a)		4.4	mg/m³	≤	10	mg/m³	yes
Carcinogenic substances, cat. 1A and 1B acc. to Regulation (EC) No. 1272/2008 (and TRGS 905) (per substance) ^{b)}	≤	0.01	mg/m³	≤	0.01	mg/m³	yes
Measurement time: 28 days after test chamber loading							
Sum VOC (C6-C16) including SVOC with LCI ^{a)}		0.30	mg/m³	≤	1.0	mg/m³	yes
Sum SVOC without LCI (C16-C22) ^{a)}		0.01	mg/m³	≤	0.1	mg/m³	yes
R-value (dimensionless)		0.05		≤	1		yes
Sum VOC without LCI	<	0.005	mg/m³	<	0.1	mg/m³	yes
Carcinogenic substances, cat. 1A and 1B acc. to Regulation (EC) No. 1272/2008 (and TRGS 905) (per substance) ^{b)}	<	0.001	mg/m³	≤	0.001	mg/m³	yes

a) For sum VOC (C6-C16) and sum SVOC (C16-C22) only substances \geq 5 $\mu g/m^3$ are considered.

b) Excluded are defined substances classified as carcinogen 1A or 1B for which a limit value can be derived for the most sensitive endpoint at which a carcinogenic potential can no longer be assumed and for which a LCI value is derived on this basis.

 $^{^1}$ If a measurement result that slightly exceeds the specification is assessed as "not fulfilled", this is based on the agreement of the "shared risk of measurement uncertainty (shared risk approach)". According to this, the probability that the statement is correct is ≥ 50 %. Similarly, a result slightly below the specification value also only has a probability of ≥ 50 % of being compliant. I.e., the risk of making a false negative statement regarding the fulfilment of the specification is just as high as the risk of making a false positive statement (more information at https://www.eco-institut.de/en/2019/07/measurement_uncertainty/).



Summary statement of conformity with AgBB 2024

The sample with the internal sample number 59983-A010, article description according to order: **SKIL REMOVE (PTKK00302 | ID13800)**, meets the emission requirements of the AgBB scheme.

Cologne, 30/07/2025

Marc-Anton Dobaj, M.Sc. Crystalline Materials

(Project management)



Laboratory report

1 Emission analysis

Test method

DIN EN 16516:2020-10 Testing and evaluation of the release of dangerous substances;

determination of emissions into indoor air

A010, Preparation of test specimen

Date: 10/06/2025

Test specimen preparation: Filled in glass dish; with a brush; application quantity 300 g/m²

drying / pre-conditioning outside of the test chamber for 72 hours

Masking of backside:

Masking of edges:

Relationship of unmasked

not applicable

not applicable

edges to surface:

Arrangement in test chamber: on tripod

Loading reference unit: area-specific [m²]

Dimensions: 2 x 25.0 cm x 20.0 cm with each 15.0 g application + 20.0 cm x 20.0 cm

with 12.0 g application

A010, Test chamber conditions according to DIN EN ISO 16000-9:2024-08

Chamber volume:

Temperature:

Relative humidity:

Air pressure:

Air:

Air change rate:

Air volority:

0.100 m³

23 °C ± 1 °C

50 % ± 5 %

normal

cleaned

0.5 h⁻¹

0.3 m (s

Air velocity: 0.3 m/s Loading: 1.4 m 2 /m 3 Specific air flow rate: 0.357 m 3 /(m 2 -h)

Starting time of the test (t0): 13/06/2025

Air sampling: 16/06/2025 (3 days after test chamber loading)

11/07/2025 (28 days after test chamber loading)



Sample A010, Volatile organic compounds after 3 days 1.1

Test objective:

Volatile organic compounds (VOC), test chamber, air sampling 3 days after test chamber loading

Method description / Analytics:

Formaldehyde and other carbonyl compounds: DIN ISO 16000-3:2023-12 (DNPH method, HPLC-DAD)

Limit of quantification: $2 \mu g/m^3$

Volatile organic compounds: DIN ISO 16000-6:2022-03 (Tenax TA®, TD-GC-MS)

Limit of quantification calibrated substances: 1 μg/m³ (1,4-Cyclohexanedimethanol, Diethylene glycol,

1,4-Butanediol: 5 μg/m³; Neodecanoic acid: 10 μg/m³)

Limit of quantification non-calibrated substances: 1 μg/m³

Test result:

Internal sample number: 59983-A010

	Substance	CAS No.	RT	Concentration+ calib. substances $\geq 1 \mu g/m^3$ uncalib. substances $\geq 1 \mu g/m^3$ DNPH $\geq 2 \mu g/m^3$	Toluene- equivalent substances ≥ 5 µg/m³	SER+	CMR Classifi- cation++	LCI AgBB 2024	R-value
			[min]	[µg/m³]	[µg/m³]	[µg/(m²·h)]		[µg/m³]	
	Aliphatic mono alcohols (n-, iso-, cyclo-) and dialcohols								
VVOC	Ethanol	64-17-5	3.48	2	< 5	0.71	III5		
	Glycols, Glycol ethers, Glycol esters								
VOC	Diethylene glycol	111-46-6	12.64	4400	1200	1600		5700	0.77
	Aldehydes								
VVOC	Formaldehyde	50-00-0		2	n. d.	0.71	Carc. 1B Muta. 2	100	0.02
	Other identified substances in addition to LCI list								
SVOC	2-n-Octyl-4-isothiazolin-3-one (OIT)	26530-20-1	28.53	8	< 5	2.9			
VOC	2-Pentanonoxime	623-40-5	10.33	4	< 5	1.4			
VOC	m/z 45 58*		10.99	2	< 5	0.71			
VOC	Phosphate compound m/z 99*		23.96	4	< 5	1.4			
SVOC	Ester m/z 115 87*		26	5	5	1.8			

⁺ identified and calibrated substances, substance specific calculated

⁺⁺ classification according to Regulation (EG) N° 1272/2008: Categories Carc. 1A, 1B and 2, Muta. 1A, 1B and 2, Repr. 1A, 1B and 2, TRGS 905: K1A, K1B, K2, M1A, M1B, M2, R1A, R1B, R2; IARC: Group 1, 2A, 2B and 3, DFG MAK-list: Categorie III1 to III5

^{*} unidentified substances, calculated as toluene equivalent reported with significant mass fragments as mass-to-charge ratio (m/z) n. d.: not determined



Carcinogenic, mutagenic, and reproductive toxic compounds*	Concentration after 3 days [µg/m³]	SERa [µg/(m² • h)]
CMR 1: VOC (incl. VVOC and SVOC) with the following categorisations: Regulation (EC) No. 1272/2008: Category Carc. 1A and 1B, Muta. 1A and 1B, Repr. 1A and 1B; TRGS 905: K1A, K1B, M1A, M1B, R1A, R1B; IARC: Group 1 and 2A; DFG (MAK list): Categories III1, III2 (sum)	< 1	< 0.36
C 1: VOC (incl. VVOC and SVOC) with the following categorisations: Regulation (EG) Nr. 1272/2008: Category Carc. 1A u. 1B; TRGS 905: K1A, K1B (sum)	<1	< 0.36

TVOC, Total volatile organic compounds	Concentration after 3 days [µg/m³]	SERa [µg/(m² • h)]
Sum of VOC according to DIN EN 16516	1200	430
Sum of VOC according to AgBB 2024	4400	1600
Sum of VOC according to eco-INSTITUT-Label	4400	1600
Sum of VOC according to DIN ISO 16000-6	1300	460

TSVOC, Total semi volatile organic compounds	Concentration after 3 days [µg/m³]	SERa [µg/(m² • h)]
Sum of SVOC according to DIN EN 16516	5	1.8
Sum of SVOC without LCI according to AgBB 2024	5	1.8
Sum of SVOC without LCI according to eco-INSTITUT-Label	13	4.6
Sum of SVOC with LCI according to AgBB 2024	< 5	< 1.8

TVVOC, Total very volatile organic compounds	Concentration after 3 days [µg/m³]	SERa [µg/(m² • h)]
Sum of VVOC according to AgBB 2024	< 5	< 1.8
Sum of VVOC according to eco-INSTITUT-Label	4	1.4

^{*}Excluding formaldehyde and acetaldehyde (Carc. 1B) due to an assumed "practical threshold" under which a significant carcinogenic risk is no longer to be expected (see Federal Institute for Risk Assessment (2006): Toxicological evaluation of formaldehyde and Federal Environment Agency (2016): Reference value for formaldehyde in indoor air and protocol of the 11th meeting of 'Ausschusses für Innenraumrichtwerte' (AIR), 11/2020). In the case of a toxicological emission assessment, a single-substance analysis of the concentrations is necessary.

In the opinion of the committee for Indoor Air Guide Values (Ausschuss für Innenraumrichtwerte) of the Federal Environment Agency, the concentration of 0.1 mg formaldehyde/m³ indoor air, based on a measurement period of half an hour, should not be exceeded, also for a short time (Bundesgesundheitsblatt 2016 · 59: 1040-1044 DOI 10.1007 / s00103 -016-2389-5 © Springer-Verlag Berlin Heidelberg 2016).



Other sums of VOC	Concentration after 3 days [µg/m³]	SERa [µg/(m² • h)]
VOC without LCI according to AgBB 2024 (sum)	< 5	< 1.8
VOC without LCI according to eco-INSTITUT-Label (sum)	10	3.6
CMR 2: VOC (incl. VVOC and SVOC) with the following categorisations: Regulation (EC) No. 1272/2008: Category Carc. 2, Muta. 2, Repr. 2; TRGS 905: K2, M2, R2; IARC: Group 2B; DFG (MAK list): Category III3 (sum)	2	0.71
Sensitising compounds with the following categorisations: DFG (MAK list): Category IV; Regulation (EC) No. 1272/2008: skin sensitising, respiratory sensitising; TRGS 907 (sum)	10	3.6
Bicyclic Terpenes (sum)	<1	< 0.36
C9 - C14 Alkanes / Isoalkanes as dekane-equivalent (sum)	< 1	< 0.36
C4 - C11 Aldehydes, acyclic, aliphatic (sum)	< 2	< 0.71
C9 - C15 Alkylated benzenes (sum)	< 1	< 0.36
Cresols (sum)	<1	< 0.36

Risk value for assessment of LCI	R-value
R-value according to eco-INSTITUT-Label	0.79
R-value according to AgBB 2024	0.77
R-value according to Belgian regulation	0.77
R-value according to EU-LCI	0.77

Note:

Due to different requirements in the respective guidelines, the calculation of TVOC, TVVOC, TSVOC and R-value may result in different values.

Short-chain carbonyl compounds (C1-C5) are quantified via HPLC acc. to DIN ISO 16000-3:2013-01. Therefore, no toluene equivalents are given for VVOC. These substances are taken into concern by means of their substance specific calibration via the sum of VVOC acc. to DIN EN 16516:2020-10. For VOC however, the substance specific calibration takes place via HPLC whereas the TVOC is calculated using the toluene equivalent determined via Tenax acc. to DIN EN 16516:2020-10.



1.2 Sample A010, Volatile organic compounds after 28 days

Test objective:

Volatile organic compounds (VOC), test chamber, air sampling 28 days after test chamber loading

Method description / Analytics:

Formaldehyde and other carbonyl compounds: DIN ISO 16000-3:2023-12 (DNPH method, HPLC-DAD)

Limit of quantification: 2 μg/m³

Volatile organic compounds: DIN ISO 16000-6:2022-03 (Tenax TA®, TD-GC-MS)

Limit of quantification calibrated substances: 1 μg/m³ (1,4-Cyclohexanedimethanol, Diethylene glycol,

1,4-Butanediol: 5 μ g/m³; Neodecanoic acid: 10 μ g/m³)

Limit of quantification non-calibrated substances: $1 \mu g/m^3$

Test result:

Internal sample number: 59983-A010

	Substance	CAS No.	RT	Concentration+ calib. substances ≥ 1 µg/m³	Toluene- equivalent	SER+	CMR Classifi- cation++	LCI AgBB 2024	R-value
				uncalib. substances ≥ 1 μg/m³	substances ≥ 5 µg/m³				
				DNPH ≥ 2 μg/m³	F / 2]	F // 2127		F (2)	
			[min]	[µg/m³]	[µg/m³]	[µg/(m²·h)]		[µg/m³]	
	Glycols, Glycol ethers, Glycol esters								
VOC	Diethylene glycol	111-46-6	12.4	300	82	110		5700	0.05
	Other identified substances in addition to LCI list								
SVOC	2-n-Octyl-4-isothiazolin-3-one (OIT)	26530-20-1	28.53	10	5	3.6			
SVOC	Ester m/z 115 87*		26	5	5	1.8			

⁺ identified and calibrated substances, substance specific calculated

⁺⁺ classification according to Regulation (EG) N° 1272/2008: Categories Carc. 1A, 1B and 2, Muta. 1A, 1B and 2, Repr. 1A, 1B and 2, TRGS 905: K1A, K1B, K2, M1A, M1B, M2, R1A, R1B, R2; IARC: Group 1, 2A, 2B and 3, DFG MAK-list: Categorie III1 to III5

 $^{^*}$ unidentified substances, calculated as toluene equivalent reported with significant mass fragments as mass-to-charge ratio (m/z)

n. d.: not determined



Carcinogenic, mutagenic, and reproductive toxic compounds*	Concentration after 28 days [µg/m³]	SERa [µg/(m² • h)]
CMR 1: VOC (incl. VVOC and SVOC) with the following categorisations: Regulation (EC) No. 1272/2008: Category Carc. 1A and 1B, Muta. 1A and 1B, Repr. 1A and 1B; IRGS 905: K1A, K1B, M1A, M1B, R1A, R1B; IARC: Group 1 and 2A; DFG (MAK list): Categories III1, III2 (sum)	<1	< 0.36
C 1: VOC (incl. VVOC and SVOC) with the following categorisations: Regulation (EG) Nr. 1272/2008: Category Carc. 1A u. 1B; TRGS 905: K1A, K1B (sum)	<1	< 0.36

TVOC, Total volatile organic compounds	Concentration after 28 days [µg/m³]	SERa [µg/(m² • h)]
Sum of VOC according to DIN EN 16516	82	29
Sum of VOC according to AgBB 2024	300	110
Sum of VOC according to eco-INSTITUT-Label	300	110
Sum of VOC according to DIN ISO 16000-6	86	31

TSVOC, Total semi volatile organic compounds	Concentration after 28 days [µg/m³]	SERa [µg/(m² • h)]
Sum of SVOC according to DIN EN 16516	10	3.6
Sum of SVOC without LCI according to AgBB 2024	10	3.6
Sum of SVOC without LCI according to eco-INSTITUT-Label	15	5.4
Sum of SVOC with LCI according to AgBB 2024	< 5	< 1.8

TVVOC, Total very volatile organic compounds	Concentration after 28 days [µg/m³]	SERa [µg/(m² • h)]
Sum of VVOC according to AgBB 2024	< 5	< 1.8
Sum of VVOC according to eco-INSTITUT-Label	<1	< 0.36

^{*}Excluding formaldehyde and acetaldehyde (Carc. 1B) due to an assumed "practical threshold" under which a significant carcinogenic risk is no longer to be expected (see Federal Institute for Risk Assessment (2006): Toxicological evaluation of formaldehyde and Federal Environment Agency (2016): Reference value for formaldehyde in indoor air and protocol of the 11th meeting of 'Ausschusses für Innenraumrichtwerte' (AIR), 11/2020). In the case of a toxicological emission assessment, a single-substance analysis of the concentrations is necessary.

In the opinion of the committee for Indoor Air Guide Values (Ausschuss für Innenraumrichtwerte) of the Federal Environment Agency, the concentration of 0.1 mg formaldehyde/m³ indoor air, based on a measurement period of half an hour, should not be exceeded, also for a short time (Bundesgesundheitsblatt 2016 · 59: 1040-1044 DOI 10.1007 / s00103 -016-2389-5 © Springer-Verlag Berlin Heidelberg 2016).



Other sums of VOC	Concentration after 28 days [µg/m³]	SERa [µg/(m² • h)]
VOC without LCI according to AgBB 2024 (sum)	< 5	< 1.8
VOC without LCI according to eco-INSTITUT-Label (sum)	<1	< 0.36
CMR 2: VOC (incl. VVOC and SVOC) with the following categorisations: Regulation (EC) No. 1272/2008: Category Carc. 2, Muta. 2, Repr. 2; TRGS 905: K2, M2, R2; IARC: Group 2B; DFG (MAK list): Category III3 (sum)	<1	< 0.36
Sensitising compounds with the following categorisations: DFG (MAK list): Category IV; Regulation (EC) No. 1272/2008: skin sensitising, respiratory sensitising; TRGS 907 (sum)	10	3.6
Bicyclic Terpenes (sum)	< 1	< 0.36
C9 - C14 Alkanes / Isoalkanes as dekane-equivalent (sum)	< 1	< 0.36
C4 - C11 Aldehydes, acyclic, aliphatic (sum)	< 2	< 0.71
C9 - C15 Alkylated benzenes (sum)	< 1	< 0.36
Cresols (sum)	<1	< 0.36

Risk value for assessment of LCI	R-value
R-value according to eco-INSTITUT-Label	0.05
R-value according to AgBB 2024	0.05
R-value according to Belgian regulation	0.05
R-value according to EU-LCI	0.05

Note:

Due to different requirements in the respective guidelines, the calculation of TVOC, TVVOC, TSVOC and R-value may result in different values.

Short-chain carbonyl compounds (C1-C5) are quantified via HPLC acc. to DIN ISO 16000-3:2013-01. Therefore, no toluene equivalents are given for VVOC. These substances are taken into concern by means of their substance specific calibration via the sum of VVOC acc. to DIN EN 16516:2020-10. For VOC however, the substance specific calibration takes place via HPLC whereas the TVOC is calculated using the toluene equivalent determined via Tenax acc. to DIN EN 16516:2020-10.

Cologne, 30/07/2025

Michael Stein, Dipl.-Chem. (Laboratory Management)



Appendix

Sampling sheet



Tel. +49 221.931245-0 / Fax +49 221.931245-33 / eco-institut.de / Geschäftsführer: Dr. Frank Kuebart, Daniel Tigges
HRB 17917 / USF-ID: DE 12265330C / Volksbank Rhein-Erft-Koln eG, IBAN: DE60370623651701900010, BIC: GENODED1FHH



List of calibrated Volatile Organic Compounds (VOC)

Aromatic hydrocarbons (30)

1,2,3-Trimethylbenzene
1,2,4-Trimethylbenzene
1,3,5-Trimethylbenzene
1-Isopropyl-2-methylbenzene
1-Isopropyl-4-methylbenzene
1,2,4,5-Tetramethylbenzene

Ethylbenzene n-Propylbenzene Isopropylbenzene (Cumene)⁴

1,3-Diisopropylbenzene
1,4-Diisopropylbenzene
n-Butylbenzene

1-Propenylbenzene (beta-Methylstyrene)

Toluene
2-Ethyltoluene
Vinyltoluene
o-Xylene
m-/p-Xylene
Styrene
Phenylacetylene

2-Phenylpropene (alpha-Methylstyrene)

4-Phenylcyclohexene
1-Phenyloctane
1-Phenyldecane²
1-Phenylundecane²
Indene
Naphthalene
1-Methylnaphthalene
2-Methylnaphthalene

Aliphatic hydrocarbons (24)

1,4-Dimethylnaphthalene

2-Methylpentane¹
3-Methylpentane¹
Methylcyclopentane
n-Pentane¹
n-Hexane

Cyclohexane Methylcyclohexane 1,4-Dimethylcyclohexane

n-Heptane

2,2,4,6,6-Pentamethylheptane

n-Octane
n-Nonane
n-Decane
n-Undecane
n-Dodecane
n-Tridecane
n-Tetradecane
n-Pentadecane
n-Hexadecane
Decahydronaphthalene

1-Octene 1-Decene 1-Dodecene 4-Vinylcyclohexene

Terpenes (12)

delta-3-Carene alpha-Pinene beta-Pinene alpha-Terpinene Longipinene Limonene Longifolene Isolongifolene beta-Caryophyllene alpha-Phellandrene Myrcene

Camphene

Fthanol

Aliphatic alcohols and ether (18)

1-Propanol¹
2-Propanol¹
2-Methyl-1-propanol
1-Butanol
tert-Butanol
1-Pentanol
1-Hexanol
Cyclohexanol
2-Ethyl-1-hexanol
1-Heptanol
1-Octanol
1-Nonanol
1-Decanol

1,4-Cyclohexandimethanol 4-Hydroxy-4-methyl-pentan-2-one

(Diacetone alcohol) Methyl-tert-butyl ether (MTBE)¹ Tetrahydrofuran (THF)

Aromatic alcohols (phenoles) (8)

Furfuryl alcohol Benzyl alcohol Phenol 2-Phenylphenol (oPP)

BHT (2,6-Di-tert-butyl-4-methylphenol)

o-Cresol m-/p-Cresol

4-Chloro-3-methylphenol (Chlorocresol)

Glycols, Glycol ether, Glycol ester (49)

Ethyleneglycol (Ethan-1,2-diol)
Propylenglycol (Propane-1,2-diol)
Diathylono glycol

Diethylene glycol Dipropylene glycol Neopentyl glycol Hexyleneglycol Ethyldiglycol

Ethylene glycol monobutyl ether Diethylene glycol methyl ether Diethylene glycol monobutyl ether Diethylene glycol phenyl ether Dipropylene glycol-dimetyl ether Dipropylene glycol mono-n-butyl ether Dipropylene glycol mono-tert-butyl ether Dipropylene glycol monomethyl ether Dipropylene glycol mono-n-propyl ether Tripropylene glycol monomethyl ether Triethylene glycol dimethyl ether 1,2-Propylene glycol dimethyl ether 1,2-Propylene glycol-n-propyl ether 1,2-Propylene glycol-n-butyl ether Butyl glycolate

2-Methylethoxyethanol
2-Propoxyethanol
2-Hexoxyethanol
2-(2-Hexoxyethoxy)ethanol
2-Phenoxyethanol
1-Methoxy-2-propanol
1-Ethoxy-2-propanol
1-tert-Butoxy-2-propanol
3-Methoxy-1-butanol
1,4-Butanediol
1,2-Dimethoxyethane
1,2-Diethoxyethane

2-Methoxyethanol

2-Ethoxyethanol

1-Methoxy-2-(2-methoxy-ethoxy)ethane

Ethylene carbonate Propylene carbonate 2-Methoxy-1-propyl acetate

Diethylene glycol monomethyl ether acetate

2-Methoxyethyl acetate 2-Ethoxyethyl acetate 2-Butoxy ethyl acetate

Dipropylene glycol monomethyl ether acetate

Propylene glycol diacetate

Texanol

TXIB (Texanol isobutyrate)

Aldehydes (26)

Formaldehyde^{1,3,4}
Acetaldehyde^{1,3,4}
Propanal^{1,3}
Butanal^{1,3}
3-Methyl-1-butanal
Pentanal
Hexanal
2-Ethylhexanal
Heptanal
Octanal
Nonanal
Decanal
Propenal (Acrolein)¹

Isobutenal (Methacrolein)³

2-Butenal 2-Pentenal³ 2-Hexenal 2-Heptenal 2-Octenal



2-Nonenal 2-Decenal 2-Undecenal Ethanedial (Glyoxal)^{1,3}

Glutaraldehyde Furfural Benzaldehyde

Ketones (14)

Acetone^{1,2} 1-Hydroxyacetone Ethylmethylketone³ Methylisobutylketone 3-Methyl-2-butanone Cyclopentanone 2-Methylcyclopentanone

Cyclohexanone

2-Methylcyclohexanone

2-Hexanone 2-Heptanone Acetophenone Isophorone

4-Methylbenzophenone²

Acids (11)

Acetic acid Propionic acid Pivalic acid Butyric acid Isobutyric acid n-Valeric acid n-Caproic acid 2-Ethylhexanoic acid n-Heptanoic acid n-Octanoic acid Neodecanoic acid

Esters and Lactones (33)

Methyl acetate1 Ethyl acetate1 Vinyl acetate¹ Propyl acetate Isopropyl acetate

2-Methoxy-1-methylethyl acetate

n-Butyl acetate Isobutylacetate 2-Ethylhexyl acetate n-Butyl formate Methyl acrylate Methyl methacrylate Butyl methacrylate Ethyl acrylate n-Butyl acrylate 2-Ethylhexyl acrylate 2-Ethylhexyl methacrylate Hexanediol diacrylate Dipropylene glycol diacrylate

Dimethyl succinate Dimethyl glutarate Dimethyl adipate Dibutyl fumarate Dibutyl maleate Diisobutyl succinate Diisobutyl glutarate Butyrolactone Dimethyl phthalate Diethyl phthalate² Dipropyl phthalate² Dibutyl phthalate² Diisobutyl phthalate²

(5-Ethyl-1,3-dioxan-5-yl)methyl acrylate

Cyclic siloxanes (5)

Hexamethylcyclotrisiloxane (D3) Octamethylcyclotetrasiloxane (D4) Decamethylcyclopentasiloxane (D5) Dodecamethylcyclohexasiloxane (D6) Tetradecamethylcycoheptasiloxane (D7)

Carcinogens (44)

Isopropylbenzene (Cumene)4

Benzene⁴ Benzophenone4

Trichloromethane (Chloroform)4

1,2-Dichloroethane4 1,2,3-Trichloropropane4 Trichloroethene4

trans-1,3-Dichloropropene4 cis-1,3-Dichloropropene4

Chloroprene4 1,3-Dichloro-2-propanol4 alpha-Chlorotoluene4

alpha,alpha,alpha-Trichlorotoluene4

1.4-Dioxane⁴ 1,2-Dibromoethane4 2-Nitropropane4 2,3-Dinitrotoluene4 2,4-Dinitrotoluene4 2,6-Dinitrotoluene4 3,4-Dinitrotoluene^{2,4} o-Anisidine4 o-Toluidine4 4-Chloro-o-toluidine4 Acrylonitrile1,4 Azobenzene^{2,4} Furan^{1,4} 2-Butanonoxime4 N-Nitrosopyrrolidine4

4-Chloroaniline4

2-Nitroanisole4 p-Cresidine4 Diethyl sulfate4 Epichlorohydrin4 1,2-Dichloropropane4 Urethane4 Acrylamide4

trans-1,4-Dichlorobut-2-ene4 1,2-Dibromo-3-chloropropane4

2-Nitrotoluene4 Quinoline4 Phenylglycidyl ether4 2,4,5-Trimethylaniline4 4-Chlorobenzotrichloride4 Nitrosodipropylamin4

Others (35)

5-Nitro-o-toluidine² 2,2'-Azobisisobutyronitrile Tetramethylsuccinonitrile

Caprolactam 2-Methylfuran 2-Pentylfuran Methenamine Diethylamine1 Triethylamine

Triethylene diamine (DABCO®)

Triethyl phosphate Tributyl phosphate²

5-Chloro-2-methyl-4-isothiazolin-3-one (CIT) 2-Methyl-4-isothiazolin-3-one (MIT) 2-n-Octyl-4-isothiazolin-3-one (OIT)

Formamide N-Methylformamide

Dimethylformamide (DMF)

Acetamide

N-Methyl-2-pyrrolidone N-Ethyl-2-pyrrolidone N-Butyl-2-pyrrolidone

Aniline5

Cyclohexyl isocyanate 5-Ethyl-1,3-dioxan-5-methanol

Dichloromethane¹ Tetrachloromethane 1,1,1-Trichloroethane 2-Chloropropane Tetrachloroethene Chlorobenzene 1,4-Dichlorobenzene 1,1-Dichlorethene1 2-Pentanone oxime

Tribromomethane (Bromoform)

2 SVOC

Analysis acc. to DIN ISO 16000-3:2023-12 (DNPH)

Carcinogens, category 1A and 1B according to Regulation (EC) No 1272/2008 and TRGS 905

When analysing with TD-GC-MS, aniline can occur as a thermal decomposition product of other substances (e.g. 1,3-Diphenylquanidine). A cold analytical method is recommended to confirm the result.

(Status: June 2025)

¹ VVOC



Definition of terms

CAS No. (Chemical Abstracts Service)

CMR

Limit of quantification (LOQ)

NIK / LCI

RT (retention time)

R value

R value according to AgBB

R-value according to Belgian regulation

R value according to eco-INSTITUT-Label

R value according to EU-LCI

SER

SVOC (semi volatile organic compound)

Toluene equivalent

TSVOC

TSVOC according to DIN EN 16516

TSVOC with LCI according to AgBB

TSVOC with LCI according to eco-INSTITUT-Label

TSVOC without LCI according to AgBB

TSVOC without LCI according to eco-INSTITUT label

TVOC

International designation standard for chemical substances

VOCs, VVOCs and SVOCs classified as carcinogenic, mutagenic or toxic for reproduction according to Regulation (EC) No. 1272/2008, TRGS 905, IARC list and DFG (MAK list)

Lower limit of quantification in the analytical method within the defined measurement uncertainty

Lowest concentration of interest; substance-specific value for health assessment of emissions from products, indicated in $\mu g/m^3$

Total time required for an analyte to pass the column (time between injection and detection of the analyte)

Sum of quotients of concentration and LCI value for all substances for which a LCI value is derived

R-value for all substances $\geq 5~\mu g/m^3$ with LCI value, calculated according to the LCI list of the AgBB scheme

R-value for all substances \geq 5 µg/m³ with LCI-value, calculated according to the LCI-list of the Belgian regulation

R-value for all substances $\geq 1~\mu g/m^3$ with LCI value, calculated according to the LCI list of the AqBB scheme

R-value for all substances \geq 5 $\mu g/m^3$ with EU-LCI value, calculated according to the EU-LCI list of the European Commission

Specific emission rate (see "Explanation of Specific Emission Rate SER")

Organic compound eluting in the retention range > C_{16} (n-hexadecane) to C_{22} (docosane)

Concentration of a substance quantified by the TIC response factor of toluene (calculation of the concentration by comparing the integral of the substance with the integral of toluene)

Sum of the concentrations of all identified and unidentified semi volatile organic compounds eluting in the retention range $> C_{16}$ (n-hexadecane) to C_{22} (docosane)

Sum of all SVOC \geq 5 µg/m³ (as toluene equivalent)

Sum of all SVOC with LCI $\geq 5 \mu g/m^3$ (quantified substance-specific)

Sum of all SVOC with LCI $\geq 1 \mu g/m^3$ (quantified substance-specific)

Sum of all SVOC without LCI \geq 5 µg/m³ (as toluene equivalent)

Sum of all calibrated SVOC without LCI \geq 1 $\mu g/m^3$ (quantified substance-specific) and all non-calibrated SVOC without LCI \geq 1 $\mu g/m^3$ (as toluene equivalent)

Sum of the concentrations of all identified and unidentified volatile organic compounds eluting in the retention range from C_6 (n-hexane) to C_{16} (n-hexadecane)



TVOC according to DIN EN 16516

TVOC according to AgBB

TVOC according to eco-INSTITUT-Label

TVOC according to ISO 16000-6

TVOC without LCI according to AgBB

TVOC without LCI according to eco-INSTITUT-Label

TVVOC

TVVOC according to AgBB

TVVOC according to eco-INSTITUT-Label

VOC (volatile organic compound)

VVOC (very volatile organic compound)

Sum of all $VOC \ge 5 \mu g/m^3$ in the retention range C_6 to C_{16} , calculated as toluene equivalent (used i.a. for M1)

Sum of all VOCs with LCI \geq 5 µg/m³ (quantified substance-specific) and all VOCs without LCI \geq 5 µg/m³ (as toluene equivalent) (used i.a. for the Blue Angel)

Sum of all calibrated VOC \geq 1 $\mu g/m^3$ (quantified substance-specific) and all non-calibrated VOC \geq 1 $\mu g/m^3$ (as toluene equivalent) (used i.a. for natureplus)

Total area of the chromatogram in the retention range C_6 – C_{16} as toluene equivalent according to DIN ISO 16000-6, Annex A.1 item 3 (used i.a. for CDPH, BIFMA and the French VOC regulation)

Sum of all VOCs without LCI $\geq 5 \mu g/m^3$ as toluene equivalent

Sum of all calibrated VOCs without LCI \geq 1 $\mu g/m^3$ (quantified substance-specific) and all non-calibrated VOCs without LCI \geq 1 $\mu g/m^3$ (as toluene equivalent)

Sum of the concentrations of all identified and unidentified very volatile organic compounds eluting in the retention range $< C_6$ (n-hexane)

Sum of all VVOC with LCI \geq 5 μ g/m³ (quantified substance-specificic) and all VVOC without LCI \geq 5 μ g/m³ (as toluene equivalent)

Sum of all calibrated VVOC \geq 1 $\mu g/m^3$ (substance-specific quantified) and all non-calibrated VVOC \geq 1 $\mu g/m^3$ (as toluene equivalent)

Organic compound eluting in the retention range from C_6 (n-hexane) to C_{16} (n-hexadecane)

Organic compound eluting in the retention range $< C_6$ (n-hexane)



Commentary on emission analysis

Test method

Measurement of the volatile organic compounds takes place in the test chamber in conditions similar to those applying in practice. Standardised test conditions are defined for the test chamber regarding loading, air exchange, relative humidity, temperature, and incoming air, based on the type of test specimen and the required guideline. These conditions and the underlying standards are to be found in the section on test methods in the laboratory report.

Air samples are taken from the test chamber at defined points in time during the continuously running test. To this end, approximately 5 L of air are collected from the test chamber at an air flow rate of 100 mL/min on Tenax and approx. 100 L at an air flow rate of 0.8 L/min on silica gel coated with DNPH (2,4-dinitrophenylhydrazine).

After thermal desorption, the substances adsorbed on Tenax are analysed using gas chromatographic separation and mass spectrometric determination. The gas chromatographic separation is performed with a slightly polar capillary column of 60 m in length.

The substances derivatised with DNPH for the determination of formaldehyde and other short-chain carbonyl compounds ($C_1 - C_6$) are analysed using high-performance liquid chromatography (HPLC).

Over 200 compounds, including volatile organic compounds ($C_6 - C_{16}$), semi-volatile organic compounds ($C_{16} - C_{22}$) and – insofar as possible with this method – also very volatile organic compounds (less than C_6) are determined and quantified individually.

All other substances – insofar as possible – are identified through comparison with a library of spectra. The quantification of these substances and non-identified substances is performed through a comparison of their signal area with the signal of toluene.

The determined substance concentrations are corrected using the recovery rate of the internal standard (toluene-d8). Identification and quantification of substances is carried out from a concentration (limit of quantification) of 1 μ g per m³ test chamber air or 2 μ g/m³ for DNPH-derivatised substances. In the case of highly loaded samples, the evaluation limit of non-calibrated substances is raised in some cases, as it is no longer possible to assign individual, small signals due to the large number of signals.

Quality assurance

The eco-INSTITUT Germany GmbH is granted flexible scope of accreditation pursuant to DIN EN ISO/IEC 17025:2018-03. The accreditation covers the analytical determination of all volatile organic compounds, including the test chamber method.

In each analysis the analytical system is checked using an external standard based on the specifications in standard DIN EN 16516:2020-10. The stability of the analytical systems is documented based on the test standard using control charts.

Laboratory performance is assessed at least once a year in inter-laboratory comparisons by comparing the results with those obtained by other laboratories for identical samples.

A blank is run prior to introducing the test specimen into the test chamber to check for the possible presence of volatile organic compounds.

The expanded measurement uncertainty U for the analytical determination of all volatile organic compounds, including the test chamber method, is estimated to 29.3 %. The calculation is based on DIN ISO 11352:2013-03 (Nordtest).



Explanation of Specific Emission Rate SER

Emission measurements are accomplished in test chambers under defined physical conditions (temperature, relative humidity, room loading, air change rate etc.).

Test chamber measurement results are directly comparable only if the investigations were accomplished under the same basic conditions.

If the differences of the physical conditions refer only to the change of air rate and/or the loading, the "SER" or "specific emission rate" can be used for comparability of the measurement results. The SER indicates how many volatile organic compounds (VOC) are released by the sample for each material unit and hour (h).

The SER can be calculated using the formula below for each proven individual component of the VOC from the data in the test report.

As material units the following are applicable:

I = unit of length (m) relation between emission and length
a = unit area (m²) relation between emission and surface
v = unit volume (m³) relation between emission and volume
u = piece unit (unit = piece) relation between emission and complete unit

From this the different dimensions for SER result:

 $\begin{array}{lll} \mbox{length-specific} & \mbox{SER}_l & \mbox{in } \mu g/(m \cdot h) \\ \mbox{surface-specific} & \mbox{SER}_a & \mbox{in } \mu g/(m^2 \cdot h) \\ \mbox{volume-specific} & \mbox{SER}_v & \mbox{in } \mu g/(m^3 \cdot h) \\ \mbox{unit-specific} & \mbox{SER}_u & \mbox{in } \mu g/(u \cdot h) \end{array}$

SER thus represents a product specific rate, which describes the mass of the volatile organic compound, which is emitted by the product per time unit at a certain time after beginning of the examination.

$$SER = q \cdot c$$

- q specific air flow rate (quotient from change of air rate and loading)
- c concentration of the measured substance(s)

The result can be indicated in milligrams (mg) in place of micro grams (μ g), whereby 1 mg = 1000 μ g.