



# Test Report

Report No.: 890653-2-AB

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Hbk/jle

Order no.: 890653

No. of appendices: 4

**Assignor:** Normann Copenhagen ApS  
Østerbrogade 70  
2100 Copenhagen  
Denmark

**Material:** HYG 2-seater sofa, with oak legs and Camira fabric, mainline flax.

**Sampling:** The test material was sampled by the assignor and received in cardboard at the Danish Technological Institute 2020-03-19.

**Method:** ANSI/BIFMA M7.1-2011 (R2016) – Standard test method for determining VOC emissions from office furniture systems, components and seating.

Details of test methods are shown in appendix 2.

**Period:** The chamber testing was carried out from 2020-03-24 to 2020-03-31.

**Result:** The estimated building concentrations ( $C_{\text{bldg}}$ ) of the tested furniture and limit of indoor air concentrations after 7 days (168 hours) are shown in the following table:

|  | Tested Seating<br>$C_{\text{bldg}}$ | Emissions Limits*<br>Seating $C_{\text{bldg}}$ |
|--|-------------------------------------|--|
| TVOC <sub>(toluene)</sub> (mg/m <sup>3</sup> ) | 0.20                                | ≤ 0.25   |
| Formaldehyde (ppb)                             | 1.3                                 | ≤ 25   |
| Total aldehydes (ppb)                          | 2.2                                 | ≤ 50   |
| 4-Phenylcyclohexene (mg/m <sup>3</sup> )       | < 0.001                             | ≤ 0.00325                                      |

\* ANSI/BIFMA M7.1 Table A1.1.

Detailed results for emissions are shown in Appendix 3.

Estimated building concentrations are shown in Appendix 4. The estimated building concentrations are below the allowable limits; hence the tested furniture is low-emitting according to LEED v4 for Interior Design and Construction, January 5, 2018.

**Note:** The test results does not apply to furniture with other materials than the tested sample, such as plywood legs or leather covers.

**Storage:** The test material will be picked up by the assignor after testing unless otherwise agreed.

**Terms:** Accredited testing was carried out in compliance with international requirements (EN/ISO/IEC 17025:2005) and in compliance with Danish Technological Institute's General Terms and Conditions regarding Commissioned Work accepted by Danish Technological Institute. The test results apply to the tested products only. This report may be quoted in extract only if the laboratory has granted its written consent.

**Date/place:** 2020-05-15, Danish Technological Institute, Taastrup, Building and Construction

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**Signature:** Test responsible

Co-signatory



**Sample information**

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Material information given by assignor:

| <b>Chain of Custody form (CoC)</b><br>ANSI BIFMA / CDPH 01350   |   |
|---|---|
| <b>Manufacturer details</b>   |   |
| Company   | Normann Copenhagen ApS  |
| Address   | Østerbrogade 70,<br>2100 Copenhagen   |
| Country   | Denmark   |
| Contact name  | Søren Stryhn-Petersen   |
| Contact title   | CTO   |
| Phone   | +45 22844421  |
| E-mail  | ssp@normann-copenhagen.com  |
| <b>Sample details</b>   |   |
| Sample ID   | HYG 2-seater sofa, with oak legs and Camira fabric, mainline flax                                 |
| Product category  | Sofas and lounge chairs   |
| Product name  | HYG & Era series – sofas and lounge chairs  |
| Manufacturers ID. no.   | [REDACTED]  |
| Date manufactured   | March 12 <sup>th</sup> , 2020   |
| Sample collection location  | [REDACTED]  |
| Samples collection time and date  | [REDACTED]  |
| Sample collected by   | [REDACTED]  |
| Number of sample pieces   | 1 pc. HYG Sofa  |
| Disposal of sample material after test:   | <input type="checkbox"/> Scrap/discard<br><input checked="" type="checkbox"/> Storage and pick-up |
| <b>Shipping details</b>   |   |
| Packed by   | Søren Stryhn-Petersen   |
| Shipping date   | March 12 <sup>th</sup> 2020   |
| Carrier   | Søren Stryhn-Petersen   |
| Air bill number   | Road delivery   |
| <b>IMPORTANT:</b><br>Please wrap samples in airtight plastic in protective cardboard box/pallet and enclose this chain of custody form.                                 |   |
| <b>Please send the sample to:</b><br>Danish Technological Institute<br>Gregersensvej, Port 3K<br>DK-2630 Taastrup<br>Attn. Helene Bendstrup Klinke – Phone + 4572202173 |   |
| Send electronic copy of test order form to: E-lab@dti.dk  |   |
| <b>Test laboratory</b>  |   |
| Reception date & initials   | 20200319/MLHON  |
| Laboratory ID   | 890653-2  |

## Sample information

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Additional information given by assignor:

# normann

C O P E N H A G E N

Appendix 1 – ANSI BIFMA/CDPH 01350

April 23<sup>rd</sup>, 2020

Test sample information:

Test sample ID: HYG sofa in full upholstery with solid oak legs

Seat material: Fabric is Camira – Mainline Flax). Rigid foam shell with steel frame, Foam cushion HR foam 60H (000023)

Leg materials: Solid oak wood, with clear lacquer

Glides: Made in HDPE

The test sample represents the following furniture with the following materials:

- 1) HYG lounge series – Sofa and lounge chairs tall and low
- 2) ERA lounge series – Sofa and lounge chairs tall and low
- 3) SUM lounge series – Sofa modules and lounge chair

The HYG series comes in following versions.

HYG sofa



HYG Lounge chairs Tall and low



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## Sample information

HYG Lounge series can be delivered with legs made of

- Solid oak wood with clear lacquer or painted
- Swivel legs in aluminium



The ERA series comes in following versions.

ERA sofas.



ERA lounge chairs:



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## Sample information

ERA sofa and lounge chairs can be delivered with legs made of:

- Plywood
- Steel with Chrome
- Steel with powder coating
- Rocking steel legs with chrome
- Aluminum swivel (only lounge chairs)

The SUM series comes as a lounge chair or sofa in modules

SUM lounge chair



SUM sofa modules



SUM lounge chair and sofa modules will be delivered with aluminum legs.

All HYG, ERA and SUM sofas and lounge chairs will be delivered as upholstery version. The upholstery version will consist of following materials:

- PUR foam shell
- Fabric: Camira - Main line flax, ERA, Oxygen, Synergy, ZAP, Oceanic, Yoredale, Zap, and Aquarius. Gabriel - Fame, Fame Hybrid, Breeze Fusion. Kvadrat - Hallingdal, Steelcut Trio and Remix.
- Leather: Sørensen leather - Ultra and Victory
- Low emission HR foam

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## Results Emissions

### Chamber testing:

The test material was unwrapped and placed in the chamber.

Photo of test material in the chamber:



|   |   |
|---|---|
| Climate chamber                             | 24 m <sup>3</sup> Polished stainless steel  |
| Temperature                                 | 23°C ± 1°C                                  |
| Relative humidity                           | 50% RH ± 5% RH                              |
| Air velocity at the surface of the specimen | 0.1 – 0.3 m/s                               |
| Air change rate (n)                         | 0.52 h <sup>-1</sup> ± 0.05 h <sup>-1</sup> |
| Area change flow rate                       | 12.4 m <sup>3</sup> /h                      |

The test material was tested in the emission chamber without prior conditioning.

Air samples were taken from the climate chamber outlet air with calibrated pumps according to ISO 16000-6 on Tenax tubes (3-4 L) and ISO 16000-3 on DNPH tubes (60 L)

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## Results Emissions

The concentration approach was applied for emission testing.

The measured chamber concentrations are shown in Tables 1 and 2.

Results from the VOC analysis are shown in Table 1.

Method: ISO 16000-6: 2011. Indoor air – Part 6: Determination of volatile organic compounds in indoor and test chamber air by active sampling on Tenax TA sorbent, thermal desorption and gas chromatography using MS/FID.

Analysis of the air sampled on Tenax was performed at the Wilhelm Klauditz Institut (WKI) under DAkkS accreditation number D-PL-11140-05-02. Report no. MAIC-2020-1268.

**Table 1: Emission of volatile organic compounds (VOCs) between n-C6 and n-C16 measured by GC-MS ( $\mu\text{g}/\text{m}^3$ )\***

| Chemical class/compound name                      | 72 hrs (3 days) |     |      |        | 168 hrs (7 days) |     |      |        |
|---|-----------------|-----|------|--------|------------------|-----|------|--------|
|   | #1              | #2  | Mean | % diff | #1               | #2  | Mean | % diff |
| <b>Aromatic hydrocarbons</b>                      |                 |     |      |        |                  |     |      |        |
| Toluene   | 2               | < 2 | 1    | >100   | 2                | < 2 | 1    | >100   |
| m,p-Xylene (m-Xylene)                             | 4               | 4   | 4    | 0      | 2                | 2   | 2    | 0      |
| Styrene   | 45              | 41  | 43   | 9      | 27               | 29  | 28   | 7      |
| o-Xylene (m-Xylene)                               | 3               | 3   | 3    | 0      | < 2              | 2   | 1    | >100   |
| Alkylbenzene (Toluene)                            | 3               | 3   | 3    | 0      | < 2              | < 2 | 0    | 0      |
| Alkylbenzene (Toluene)                            | 7               | 6   | 7    | 15     | 3                | 4   | 4    | 29     |
| Alkylbenzene (Toluene)                            | 11              | 10  | 11   | 10     | 4                | 5   | 5    | 22     |
| Alkylbenzene (Toluene)                            | 12              | 12  | 12   | 0      | 6                | 7   | 7    | 15     |
| Alkylbenzene (Toluene)                            | 5               | 4   | 5    | 22     | 3                | 3   | 3    | 0      |
| <b>Aliphatic hydrocarbons</b>                     |                 |     |      |        |                  |     |      |        |
| Cyclohexane                                       | 9               | 8   | 9    | 12     | 5                | 5   | 5    | 0      |
| saturated aliphatic hydrocarbons C9-C16 (Toluene) | 100             | 91  | 96   | 9      | 55               | 58  | 57   | 5      |
| saturated aliphatic hydrocarbons C9-C16 (Toluene) | 18              | 16  | 17   | 12     | 9                | 10  | 10   | 11     |
| saturated aliphatic hydrocarbons C9-C16 (Toluene) | 5               | 5   | 5    | 0      | 3                | 3   | 3    | 0      |
| saturated aliphatic hydrocarbons C9-C16 (Toluene) | 4               | 4   | 4    | 0      | 2                | 3   | 3    | 40     |
| saturated aliphatic hydrocarbons C9-C16 (Toluene) | 6               | 5   | 6    | 18     | 4                | 4   | 4    | 0      |
| saturated aliphatic hydrocarbons C9-C16 (Toluene) | 2               | < 2 | 1    | >100   | < 2              | < 2 | 0    | 0      |
| <b>Cycloalkanes</b>                               | < 2             | < 2 | < 2  | 0      | < 2              | < 2 | < 2  | 0      |
| <b>Terpenes</b>                                   |                 |     |      |        |                  |     |      |        |
| alpha-Pinene                                      | 2               | 2   | 2    | 0      | < 2              | < 2 | 0    | 0      |
| Terpene (Toluene)                                 | 5               | 4   | 5    | 22     | 4                | 5   | 5    | 22     |
| Terpene (Toluene)                                 | 5               | 5   | 5    | 0      | 3                | 4   | 4    | 29     |
| Longifolene                                       | 5               | 5   | 5    | 0      | 3                | 4   | 4    | 29     |
| <b>Alcohols</b>                                   |                 |     |      |        |                  |     |      |        |
| Diacetone alcohol                                 | 120             | 115 | 118  | 4      | 50               | 60  | 55   | 18     |
| <b>Glycols/Glycol ethers</b>                      |                 |     |      |        |                  |     |      |        |
|   | < 2             | < 2 | < 2  | 0      | < 2              | < 2 | < 2  | 0      |
| <b>Aldehydes</b>                                  |                 |     |      |        |                  |     |      |        |
| n-Hexanal   | 2               | 2   | 2    | 0      | < 2              | < 2 | 0    | 0      |
| Benzaldehyde                                      | 4               | 3   | 4    | 29     | 2                | 2   | 2    | 0      |

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## Results Emissions

| Chemical class/compound name                   | 72 hrs (3 days) |     |      |        | 168 hrs (7 days) |     |      |        |
|--|-----------------|-----|------|--------|------------------|-----|------|--------|
|  | #1              | #2  | Mean | % diff | #1               | #2  | Mean | % diff |
| <b>Ketones</b>                                 |                 |     |      |        |                  |     |      |        |
| 2-Pentanone, 4-hydroxy- (Toluene)              | 2               | 2   | 2    | 0      | < 2              | < 2 | 0    | 0      |
| <b>Halocarbons</b>                             |                 |     |      |        |                  |     |      |        |
| Chloroprene derivatives (Toluene)              | 4               | 4   | 4    | 0      | 3                | 3   | 3    | 0      |
| Chloroprene derivatives (Toluene)              | 20              | 18  | 19   | 11     | 12               | 13  | 13   | 8      |
| Chloroprene derivatives (Toluene)              | 14              | 12  | 13   | 15     | 9                | 10  | 10   | 11     |
| <b>Acids</b>                                   |                 |     |      |        |                  |     |      |        |
| Acetic acid                                    | 12              | < 2 | 6    | >100   | 2                | 3   | 3    | 40     |
| 2-Ethylhexanoic acid                           | 10              | 10  | 10   | 0      | 7                | 9   | 8    | 25     |
| <b>Esters</b>                                  |                 |     |      |        |                  |     |      |        |
| Ethyl acetate                                  | 12              | 12  | 12   | 0      | 7                | 8   | 8    | 13     |
| 1-Methoxy-2-propyl acetate                     | 2               | < 2 | 1    | >100   | < 2              | < 2 | 0    | 0      |
| Benzoic acid, 4-ethoxy-, ethyl ester (Toluene) | 8               | 7   | 8    | 13     | 6                | 6   | 6    | 0      |
| <b>Others</b>                                  |                 |     |      |        |                  |     |      |        |
| Octamethylcyclotetrasiloxane (D4)              | 6               | 6   | 6    | 0      | 3                | 4   | 4    | 29     |
| N,N-Dimethylbenzylamine                        | 17              | 15  | 16   | 13     | 9                | 12  | 11   | 29     |
| Decamethylcyclopentasiloxane (D5)              | 3               | 3   | 3    | 0      | 2                | 2   | 2    | 0      |
| 4-tert-Butylphenol (Toluene)                   | 9               | 8   | 9    | 12     | 8                | 9   | 9    | 12     |
| 2,4,6-Tri-tert-butylphenol (Toluene)           | 15              | 12  | 14   | 22     | 12               | 12  | 12   | 0      |
| <b>Sums</b>                                    |                 |     |      |        |                  |     |      |        |
| TVOC Toluene (ISO 16000-6)                     | 949             | 849 | 899  | 11     | 603              | 650 | 627  | 8      |
| Sum VOC (C6-C16)                               | 513             | 457 | 485  | 12     | 267              | 301 | 284  | 12     |

- \* Single substances/volatile compounds were quantified with pure reference standards, and in some cases the substances shown in subscript were used for the quantification.
- < 1 Not detected (< 1 µg/m<sup>3</sup>)  
Measured concentrations just above limit of quantification (LOQ) of 1 µg/ m<sup>2</sup>h will result in higher standard deviation from mean value.

Definitions according to ISO 16000-6:

- VOC (C6-C16): Volatile organic compounds, between hexane (C6) and hexadecane (C16)  
 VVOC (<C6): Very volatile organic compounds, eluting before hexane, not included in TVOC  
 SVOC (>C16): Semi-volatile organic compounds, eluting after hexadecane, not included in TVOC  
 TVOC: Total volatile organic compounds is the sum of all VOCs eluting between C6 and C16, quantified as toluene equivalents.

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## Results Emissions

Results from aldehyde analysis are shown in Table 2.

Method: ISO 16000-3: 2011. Indoor Air – Part 3: Determination of formaldehyde and other carbonyl compounds – Active sampling method.

Analysis of the air sampled on DNPH was performed at the Danish Technological Institute under DANAK accreditation number 90. Report no. 923787.

**Table 2: Emission of lower aldehydes (C1-C4) by HPLC analysis ( $\mu\text{g}/\text{m}^3$ )**

| Compound name | 72 hrs (3 days) |     |      |        | 168 hrs (7 days) |     |      |        |
|---------------|-----------------|-----|------|--------|------------------|-----|------|--------|
|               | #1              | #2  | Mean | % diff | #1               | #2  | Mean | % diff |
| Formaldehyde  | < 2             | < 2 | < 2  | 0      | < 2              | < 2 | < 2  | 0      |
| Acetaldehyde  | < 2             | < 2 | < 2  | 0      | < 2              | < 2 | < 2  | 0      |
| Propanal      | -               | < 2 | < 2  | > 100  | -                | -   | -    | 0      |
| Butanal       | -               | -   | -    | 0      | -                | -   | -    | 0      |
| Arolein       | -               | -   | -    | 0      | -                | -   | -    | 0      |

- Not detected. Limit of detection (LOD) is  $0.5 \mu\text{g}/\text{m}^3$  (formaldehyde, acetaldehyde, butanal),  $0.8 \mu\text{g}/\text{m}^3$  (propanal) and  $3.3 \mu\text{g}/\text{m}^3$  (acrolein).  
Measured concentrations just above limit of quantification (LOQ) of  $1 \mu\text{g}/\text{m}^2\text{h}$  will result in higher standard deviation from mean value

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## Building Concentration

### Estimated indoor air concentrations in a building

The tested furniture is seating, hence the measured concentrations from chamber testing must be adjusted to a standard office scenario using the default ventilation rate of per seating unit:  $Q=24.8 \text{ m}^3/\text{h}$ .

The concentration approach was applied for emission testing of the 2-seater sofa. Emissions correspond to 2 chairs (1-seaters), however the emissions were considered as "1 chair" in order to reflect worst case. Hence, calculations of building concentrations were carried out reflecting the worst-case emission scenario of seating in a building.

According to Table 11.1 (ANSI/BIFMA M7.1) the standard office environment parameter "modelled air flow (Q)" for seating is  $24.8 \text{ m}^3/\text{h}$  for both open plan working stations and private office work stations.

The chamber ventilation rate was  $Q_{\text{chamber}} = 12.4 \text{ m}^3/\text{h}$ , hence the measured chamber concentrations ( $C_{\text{chamber}}$ ) in Tables 1 and 2 must be multiplied by a factor:

$$Q_{\text{chamber}} = 12.4 \text{ m}^3/\text{h}$$

$$Q_{\text{bdg}} = 24.8 \text{ m}^3/\text{h}$$

$$C_{\text{bdg}} \times Q_{\text{bdg}} = C_{\text{chamber}} \times Q_{\text{chamber}}$$

$$C_{\text{bdg}} = \frac{12.4 \text{ m}^3/\text{h}}{24.8 \text{ m}^3/\text{h}} \times C_{\text{chamber}} = 0.50 \times C_{\text{chamber}}$$

Estimated building concentrations are shown in Tables 4 and 5.

The building concentrations after 14 days (336 h) were calculated from the emission factors after 3 days (72 h) and 7 days (168 h) by the power-of-law model according to the formulas given in ANSI/BIFMA M7.1 chapter 11.2.

The estimated building concentrations were calculated and results are shown in Tables 3 and 4.

**Table 3: Estimated building concentrations  $C_{\text{bdg}}$  of volatile organic compounds (VOCs) between n-C6 and n-C16 measured by GC-MS ( $\mu\text{g}/\text{m}^3$ )\***

| Chemical class/compound name                      | 3 days | 7 days | 14 days |
|---|--------|--------|---------|
| <b>Aromatic hydrocarbons</b>                      | 0.5    | 0.5    | 0.5     |
| Toluene   | 2.0    | 1.0    | 0.6     |
| m,p-Xylene (m-Xylene)                             | 21.5   | 14.0   | 9.9     |
| Styrene   | 1.5    | 0.5    | 0.2     |
| o-Xylene (m-Xylene)                               | 1.5    | < 0.1  | < 0.1   |
| Alkylbenzene (Toluene)                            | 3.3    | 1.8    | 1.1     |
| Alkylbenzene (Toluene)                            | 5.3    | 2.3    | 1.1     |
| Alkylbenzene (Toluene)                            | 6.0    | 3.3    | 2.0     |
| Alkylbenzene (Toluene)                            | 2.3    | 1.5    | 1.1     |
| Alkylbenzene (Toluene)                            |        |        |         |
| <b>Aliphatic hydrocarbons</b>                     | 4.3    | 2.5    | 1.6     |
| Cyclohexane                                       | 47.8   | 28.3   | 18.4    |
| saturated aliphatic hydrocarbons C9-C16 (Toluene) | 8.5    | 4.8    | 3.0     |
| saturated aliphatic hydrocarbons C9-C16 (Toluene) | 2.5    | 1.5    | 1.0     |
| saturated aliphatic hydrocarbons C9-C16 (Toluene) | 2.0    | 1.3    | 0.9     |
| saturated aliphatic hydrocarbons C9-C16 (Toluene) | 2.8    | 2.0    | 1.5     |
| saturated aliphatic hydrocarbons C9-C16 (Toluene) | 0.5    | < 0.1  | < 0.1   |
| saturated aliphatic hydrocarbons C9-C16 (Toluene) | -      | -      | -       |
| <b>Cycloalkanes</b>                               |        |        |         |
| <b>Terpenes</b>                                   | 1.0    | < 0.1  | < 0.1   |

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## Building Concentration

|  |       |       |       |
|--|-------|-------|-------|
| alpha-Pinene                                   | 2.3   | 2.3   | 2.3   |
| Terpene (Toluene)                              | 2.5   | 1.8   | 1.3   |
| Terpene (Toluene)                              | 2.5   | 1.8   | 1.3   |
| Longifolene                                    |       |       |       |
| <b>Alcohols</b>                                | 58.8  | 27.5  | 14.8  |
| Diacetone alcohol                              | -     | -     | -     |
| <b>Glycols/Glycol ethers</b>                   |       |       |       |
| <b>Aldehydes</b>                               | 1.0   | < 0.1 | < 0.1 |
| n-Hexanal                                      | 1.8   | 1.0   | 0.6   |
| Benzaldehyde                                   |       |       |       |
| <b>Ketones</b>                                 | 1.0   | < 0.1 | < 0.1 |
| 2-Pentanone, 4-hydroxy- (Toluene)              |       |       |       |
| <b>Halocarbons</b>                             | 2.0   | 1.5   | 1.2   |
| Chloroprene derivatives (Toluene)              | 9.5   | 6.3   | 4.4   |
| Chloroprene derivatives (Toluene)              | 6.5   | 4.8   | 3.7   |
| Chloroprene derivatives (Toluene)              |       |       |       |
| <b>Acids</b>                                   | 3.0   | 1.3   | 0.6   |
| Acetic acid                                    | 5.0   | 4.0   | 3.3   |
| 2-Ethylhexanoic acid                           |       |       |       |
| <b>Esters</b>                                  | 6.0   | 3.8   | 2.6   |
| Ethyl acetate                                  | 0.5   | < 0.1 | < 0.1 |
| 1-Methoxy-2-propyl acetate                     | 3.8   | 3.0   | 2.5   |
| Benzoic acid, 4-ethoxy-, ethyl ester (Toluene) |       |       |       |
| <b>Others</b>                                  | 3.0   | 1.8   | 1.1   |
| Octamethylcyclotetrasiloxane (D4)              | 8.0   | 5.3   | 3.7   |
| N,N-Dimethylbenzylamine                        | 1.5   | 1.0   | 0.7   |
| Decamethylcyclopentasiloxane (D5)              | 4.3   | 4.3   | 4.3   |
| 4-tert-Butylphenol (Toluene)                   | 6.8   | 6.0   | 6.0   |
| 2,4,6-Tri-tert-butylphenol (Toluene)           |       |       |       |
| <b>Sums</b>                                    | 332.0 | 195.8 | 127.1 |
| TVOC Toluene (ISO 16000-6)                     | 242.5 | 142.0 | 91.7  |

**Table 4: Estimated building concentrations  $C_{\text{bdg}}$  of lower aldehydes (C1-C4) by HPLC analysis (ppb)**

| Compound name | 3 days | 7 days | 14 days |
|---------------|--------|--------|---------|
| Formaldehyde  | 1.6    | 1.3    | 1.2     |
| Acetaldehyde  | 0.8    | 0.7    | 0.6     |
| Propanal      | < 0.1  | -      | -       |
| Butanal       | -      | -      | -       |
| Acrolein      | -      | -      | -       |

According to ANSI/BIFMA M7.1 Table A1.1 and ANSI/BIFMA X7.1 Standard for Formaldehyde and TVOC Emissions from Low-emitting Office Furniture and Seating Table A1.1, the evaluation of indoor air concentrations from furniture in a building must meet the criteria limits for giving LEED EQ Credit 4.5 Low-Emitting Materials, Systems Furniture and Seating. The results are shown in Table 5.

Higher aldehydes analysed by GC-MS VOC analysis (Table 3) were converted to ppm and included in the sum of total aldehydes (Table 5).

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## Building Concentration

**Table 5: Estimated building concentrations of tested furniture and limit of indoor air concentrations after 7 days.**

|  | Tested Seating<br>$C_{\text{bldg}}$ | Emissions Limits*<br>Seating $C_{\text{bldg}}$ |
|--|-------------------------------------|--|
| TVOC <sub>(toluene)</sub> (mg/m <sup>3</sup> ) | 0.20                                | ≤ 0.25   |
| Formaldehyde (ppb)                             | 1.3                                 | ≤ 25   |
| Total aldehydes (ppb)                          | 2.2                                 | ≤ 50   |
| 4-Phenylcyclohexene (mg/m <sup>3</sup> )       | < 0.001                             | ≤ 0.00325                                      |

\* ANSI/BIFMA M7.1 Table A1.1.

The building concentrations are below the maximum allowable limit after 7 days, and hence is low emitting furniture and compliant with LEED v4 for interior design and construction.

The estimated building concentrations according to the above ANSI/BIFMA office furniture calculations of volatile compounds in air after 3, 7 and 14 days, respectively are shown Table 6 with reference to CDPH 01350 table 4-1, where the allowable concentration limit CREL (Chronic Reference Exposure Level) are stated for evaluation of the 14 days (336 hrs) building concentrations.

Note: CDPH01350 evaluation is only relevant for classroom furniture according to LEED v4.

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## Building Concentration

**Table 6: Estimated building concentrations of tested furniture at 3, 7 and 14 days of target CREL VOCs from Proposition 65 and Table 4-1 in CDPH 01350\*.**

| Compound Name                            | CAS No.     | C <sub>(3 d)</sub><br>(µg/m <sup>3</sup> ) | C <sub>(7 d)</sub><br>(µg/m <sup>3</sup> ) | C <sub>(14 d)</sub><br>(µg/m <sup>3</sup> ) | Allowable C <sub>(14 d)</sub> **<br>(µg/m <sup>3</sup> ) |
|--|-------------|--|--|---|--|
| Acetaldehyde                             | 000075-07-0 | 1.5  | 1.2  | 1.0   | 70   |
| Benzene                                  | 000071-43-2 | -  | -  | -   | 1.5  |
| Carbon disulfide                         | 000075-15-0 | -  | -  | -   | 400  |
| Carbon tetrachloride                     | 000056-23-5 | -  | -  | -   | 20   |
| Chlorobenzene                            | 000108-90-7 | -  | -  | -   | 500  |
| Chloroform                               | 000067-66-3 | -  | -  | -   | 150  |
| Dichlorobenzene (1.4-)                   | 000106-46-7 | -  | -  | -   | 400  |
| Dichloroethylene (1.1)                   | 000075-35-4 | -  | -  | -   | 35   |
| Dimethylformamide (N.N-)                 | 000068-12-2 | -  | -  | -   | 40   |
| Dioxane (1.4-)                           | 000123-91-1 | -  | -  | -   | 1500   |
| Epichlorohydrin                          | 000106-89-8 | -  | -  | -   | 1.5  |
| Ethylbenzene                             | 000100-41-4 | -  | -  | -   | 1000   |
| Ethylene glycol                          | 000107-21-1 | -  | -  | -   | 200  |
| Ethylene glycol monoethyl ether          | 000110-80-5 | -  | -  | -   | 35   |
| Ethylene glycol monoethyl ether acetate  | 000111-15-9 | -  | -  | -   | 150  |
| Ethylene glycol monomethyl ether         | 000109-86-4 | -  | -  | -   | 30   |
| Ethylene glycol monomethyl ether acetate | 000110-49-6 | -  | -  | -   | 45   |
| Formaldehyde                             | 000050-00-0 | 2.0  | 1.7  | 1.4   | 9  |
| Hexane (n-)                              | 000110-54-3 | -  | -  | -   | 3500   |
| Isophorone                               | 000078-59-1 | -  | -  | -   | 1000   |
| Isopropanol                              | 000067-63-0 | -  | -  | -   | 3500   |
| Methyl chloroform                        | 000071-55-6 | -  | -  | -   | 500  |
| Methylene chloride                       | 001634-04-4 | -  | -  | -   | 200  |
| Methyl <i>t</i> -butyl ether             | 000075-09-2 | -  | -  | -   | 4000   |
| Naphthalene                              | 000091-20-3 | -  | -  | -   | 4.5  |
| Phenol                                   | 000108-95-2 | -  | -  | -   | 100  |
| Propylene glycol monomethyl ether        | 000107-98-2 | -  | -  | -   | 3500   |
| Styrene                                  | 000100-42-5 | 21.5                                       | 14.0                                       | 9.9   | 450  |
| Tetrachloroethylene                      | 000127-18-4 | -  | -  | -   | 17.5   |
| Toluene                                  | 000108-88-3 | 0.5  | 0.5  | 0.5   | 150  |
| Trichloroethylene                        | 000079-01-6 | -  | -  | -   | 300  |
| Vinyl acetate                            | 000108-05-4 | -  | -  | -   | 100  |
| Xylenes, technical mixture (o.m.p)       | 000106-42-3 | 3.5  | 1.5  | 0.8   | 350  |
|  | 000108-38-3 |  |  |   |  |
|  | 000095-47-6 |  |  |   |  |
|  | 001330-20-7 |  |  |   |  |
| 179601-23-1                              |             |  |  |   |  |

- Not detected

\*CDPH 01350 (2017) Standard method for the testing and evaluation of volatile organic chemical emissions from indoor sources using environmental chambers. Version 1.2.

\*\*All maximum allowable concentrations are one-half the corresponding CREL adopted by Office of Environmental Health Hazard Assessment (OEHHA) agency of California Environmental Protection Agency (CalEPA).

Results from Table 6 show the building concentrations of target CREL VOCs are below the maximum allowable limit after 14 days.