



## **Dräger X-pid<sup>®</sup> 8500/9000/9500** Product Specification V9 – 13.09.2022



Dräger. Technology for Life®

## **1** Overview sensor unit

The technical data of the control unit has to be considered separately in its instructions of use.

| Ambient conditions:             |   |  |  |  |  |
|---------------------------------|---|--|--|--|--|
| at operation                    | –10 to +35 °C   |  |  |  |  |
|                                 | 700 to 1300 hPa   |  |  |  |  |
|                                 | 10 to 95 % RH   |  |  |  |  |
|                                 |   |  |  |  |  |
| Protection class                | IP 54   |  |  |  |  |
|                                 | Gas inlets and outlets have to be protected from water and dust. The water and dust filter has to be fitted at all times. |  |  |  |  |
|                                 |   |  |  |  |  |
| Operating times                 | typically 8 h,  |  |  |  |  |
|                                 | reduces with lower ambient temperatures   |  |  |  |  |
| Dimensions                      | ca. 132 x 281 x 56 mm (W x H x T)   |  |  |  |  |
| Weight                          | ca. 880 g   |  |  |  |  |
| Approvals:                      |   |  |  |  |  |
| ATEX                            | II 1G Ex ia IIC T4 Ga   |  |  |  |  |
| IECEx                           | Ex ia IIC T4 Ga   |  |  |  |  |
| cCSAus                          | Class I, Div. 1 Group A, B, C & D T4, Ex ia   |  |  |  |  |
|                                 | Class I, Zone 0, A/Ex ia IIC T4 Ga  |  |  |  |  |
| CE Marking                      | RED (Directive 2014/53/EU)  |  |  |  |  |
| _                               | ATEX (Directive 2014/34/EU)   |  |  |  |  |
| Measurement mode Se             | eeker: (only 9x00)  |  |  |  |  |
| Sensor                          | 10.6 eV PID (Seeker-PID)  |  |  |  |  |
|                                 | Sensitive for compounds < 10.6 eV ionization energy   |  |  |  |  |
| Precision <sup>1</sup>          | < 2 % at 10.0 ppm isobutylene   |  |  |  |  |
| ( <i>k</i> = 1, ~68 %)          | < 2 % at 5.00 ppm benzene   |  |  |  |  |
| Precision <sup>1</sup>          | < 4 % at 10.0 ppm isobutylene   |  |  |  |  |
| ( <i>k</i> = 2, ~95 %)          | < 4 % at 5.00 ppm benzene   |  |  |  |  |
| Limit of detection <sup>2</sup> | 0.01 ppm (isobutylene response)   |  |  |  |  |
| Upper range <sup>3</sup>        | 60.0 ppm (isobutylene response)   |  |  |  |  |
|                                 |   |  |  |  |  |

<sup>&</sup>lt;sup>1</sup> Measure for the repeatability of measurement results at identical circumstances. The precision was determined as the *k*-fold relative standard deviation (standard or expanded uncertainty, ~68 or ~95 % confidence interval) for measurements over the course of two days of five hours each with a break but no calibration between days. Therefore, with sufficient sample size, ~68% or 95 % of measurement results will vary less than the corresponding precision. The precision does not describe the difference from measurement results and true concentration that may actually be present. <sup>2</sup> Lowest concentration that can be detected thus at which the sensor (PID) returns a signal. The limit of detection (LOD)

<sup>&</sup>lt;sup>2</sup> Lowest concentration that can be detected thus at which the sensor (PID) returns a signal. The limit of detection (LOD) depends on the sensitivity of the sensor. The LOD is valid for 100 % sensitivity of the respective PID, a parameter which is determined during calibration and that can be viewed in the archive in the user interface of the control unit.

<sup>&</sup>lt;sup>3</sup> Highest concentration that can be measured thus at which the sensor (PID) returns a signal within the specification of the device.

| Resolution                                       | 0.01 ppm from 0 to 9.99 ppm<br>0.1 ppm from 10 to 99.9 ppm  |  |  |  |  |  |
|--|---|--|--|--|--|--|
|  | 1 ppm from 100 ppm  |  |  |  |  |  |
| Measurement duration                             | direct reading  |  |  |  |  |  |
| Response time t <sub>90</sub>                    | ca. 45 seconds (isobutylene, without hose)  |  |  |  |  |  |
| Selectivity                                      | Sum concentration assuming isobutylene response for e<br>signal.<br>No selectivity between detectable compounds.  |  |  |  |  |  |
| Measurement mode Anal                            | lysis:  |  |  |  |  |  |
| Sensor   | 10.6 eV PID (Analysis-PID) after separation with gas<br>chromatography<br>Sensitive for compounds < 10.6 eV ionization energy and   |  |  |  |  |  |
| Due sisie u1                                     | boiling point < 150 °C  |  |  |  |  |  |
| Precision <sup>1</sup><br>( <i>k</i> = 1, ~68 %) | < 2 % at 10.0 ppm isobutylene<br>< 2 % at 5.00 ppm benzene  |  |  |  |  |  |
| $(K = 1, \sim 00 \%)$<br>Precision <sup>1</sup>  | < 2 % at 5.00 ppm benzene<br>< 4 % at 10.0 ppm isobutylene  |  |  |  |  |  |
| (k = 2, ~95%)                                    | < 4 % at 10.0 ppm isobutylene<br>< 4 % at 5.00 ppm benzene  |  |  |  |  |  |
| Limit of detection <sup>2</sup>                  | Compound-specific, see technical manual<br>0.07 ppm isobutylene<br>0.02 ppm benzene   |  |  |  |  |  |
| Limit of quantification <sup>4</sup>             | Compound-specific, see technical manual<br>0.20 ppm isobutylene<br>0.05 ppm benzene   |  |  |  |  |  |
| Upper range <sup>3</sup>                         | Compound-specific, see technical manual<br>100 ppm isobutylene<br>25.0 ppm benzene  |  |  |  |  |  |
| Resolution                                       | 0.01 ppm from 0 to 9.99 ppm<br>0.1 ppm from 10 to 99.9 ppm<br>1 ppm from 100 ppm  |  |  |  |  |  |
| Analysis duration                                | Compound-specific, limited by least volatile compound<br>20 s isobutylene analysis program<br>30 s benzene analysis program<br>30 s isobutylene & benzene analysis program  |  |  |  |  |  |
| Response time t <sub>90</sub>                    | none<br>(if sample concentration at start of the analysis is at device)   |  |  |  |  |  |
| Selectivity                                      | Compound-specific, see technical manual<br>For benzene there are no cross-sensitivities to toluene,<br>ethylbenzene, xylene isomers, n-hexane and many other VOC<br>with different volatility. Benzene has a known cross-sensitivity<br>to cyclohexane. |  |  |  |  |  |

<sup>&</sup>lt;sup>4</sup> Lowest concentration that can be measured thus at which the sensor (PID) returns a signal within the specification, i.e. the precision of the device. The limit of quantification (LOQ) depends on the sensitivity of the sensor. The LOQ is valid for 100 % sensitivity of the Analysis-PID, a parameter which is determined during calibration and that can be viewed in the archive in the user interface of the control unit.

The LOQ is equal to three-times the LOD. At lower concentrations rising signals ("peaks") in the chromatogram of measurement mode Analysis do not sufficiently differ from the noise of the sensor.

## 2 Target compounds (Analysis)

In measurement mode Analysis, the Dräger X-pid<sup>®</sup> x000 is limited to the target compounds benzene and 1,3-butadiene. For the Dräger X-pid<sup>®</sup> x500 the following target compounds are qualified and quantified.

| Target compounds           | CAS number | <i>tℝ</i> <sup>5</sup> , s | LOD <sup>2</sup> , ppm | LOQ⁴,ppm | Upper range <sup>3</sup> , ppm |
|----------------------------|------------|----------------------------|------------------------|----------|--------------------------------|
| Acetone                    | 67-64-1    | 8.1                        | 0.17                   | 0.50     | 50.0                           |
| Acrolein                   | 107-02-8   | 7.8                        | 0.33                   | 1.00     | 100                            |
| Benzene                    | 71-43-2    | 19.3                       | 0.02                   | 0.05     | 25.0                           |
| Butadiene, 1,3-            | 106-99-0   | 6.4                        | 0.07                   | 0.20     | 25.0                           |
| Butyl acetate              | 123-86-4   | 64.3                       | 0.67                   | 2.00     | 220                            |
| Butyraldehyde              | 123-72-8   | 12.2                       | 4.00                   | 12.0     | 210                            |
| Carbon disulfide           | 75-15-0    | 9.8                        | 0.33                   | 1.00     | 110                            |
| Cyclohexane                | 110-82-7   | 20.3                       | 0.67                   | 2.00     | 200                            |
| Dichloroethene, 1,1-       | 75-35-4    | 8.9                        | 0.07                   | 0.20     | 50.0                           |
| Dichloroethene, cis-1,2-   | 156-59-2   | 13.4                       | 0.07                   | 0.20     | 50.0                           |
| Dichloroethene, trans-1,2- | 156-60-5   | 10.9                       | 0.07                   | 0.20     | 50.0                           |
| Diethylcarbonat            | 105-58-8   | 48.4                       | 4.33                   | 13.0     | 1200                           |
| Dimethylcarbonate          | 616-38-6   | 14.6                       | 1.67                   | 5.00     | 500                            |
| Ethanol                    | 64-17-5    | 7.5                        | 1.67                   | 5.00     | 500                            |
| Ethylbenzene               | 100-41-4   | 88.7                       | 1.00                   | 3.00     | 300                            |
| Ethylene                   | 74-85-1    | 5.37                       | 0.15                   | 0.45     | 100                            |
| Ethylene oxide             | 75-21-8    | 6.8                        | 0.33                   | 1.00     | 100                            |
| Ethylmethylcarbonat        | 623-53-0   | 25.5                       | 4.17                   | 12.5     | 2000                           |
| Heptane, n-                | 142-82-5   | 27.1                       | 5.00                   | 15.0     | 500                            |
| Hexane, n-                 | 110-54-3   | 13.7                       | 0.33                   | 1.00     | 100                            |
| Hydrogen sufide            | 7783-06-4  | 5.5                        | 0.17                   | 0.50     | 55.0                           |
| Isobutylene                | 115-11-7   | 6.3                        | 0.07                   | 0.20     | 100                            |
| Isopropyl alcohol          | 67-63-0    | 9.1                        | 1.00                   | 3.00     | 200                            |
| Methyl acrylate            | 96-33-3    | 14.4                       | 0.67                   | 2.00     | 200                            |
| Methyl bromide             | 74-83-9    | 6.8                        | 0.17                   | 0.50     | 100                            |
| Methyl methacrylate        | 80-62-6    | 27.7                       | 2.50                   | 7.50     | 275                            |
| Methylbutyraldehyde, 2-    | 96-17-3    | 19.3                       | 0.33                   | 1.00     | 110                            |
| Phosphine                  | 7803-51-2  | 5.3                        | 0.67                   | 2.00     | 100                            |
| Propylene oxide            | 75-56-9    | 8.2                        | 0.17                   | 0.50     | 25.0                           |
| Propanol, 1-               | 71-23-8    | 11.6                       | 5.00                   | 15.0     | 550                            |
| Styrene                    | 100-42-5   | 111.3                      | 1.00                   | 3.00     | 300                            |
| Tetrachloroethylene        | 127-18-4   | 58.9                       | 0.67                   | 2.00     | 150                            |
| Tetrahydrofuran            | 109-99-9   | 16.5                       | 1.00                   | 3.00     | 200                            |
| Toluene                    | 108-88-3   | 41.6                       | 0.33                   | 1.00     | 100                            |
| Trichloroethylene          | 79-01-6    | 24.9                       | 0.33                   | 1.00     | 100                            |
| Vinyl acetate              | 108-05-4   | 11.9                       | 0.33                   | 1.00     | 55.0                           |
| Vinyl chloride             | 75-01-4    | 6.3                        | 0.33                   | 1.00     | 100                            |
| Xylene, m-                 | 108-38-3   | 95.7                       | 1.00                   | 3.00     | 300                            |
| Xylene, o-                 | 95-47-6    | 114.5                      | 1.00                   | 3.00     | 300                            |
| Xylene, p-                 | 106-42-3   | 96.6                       | 1.00                   | 3.00     | 300                            |

## 2.1 Qualified compounds

For the Dräger X-pid<sup>®</sup> x500, further target compounds are qualified, but not quantified. Not always was the measurement range experimentally determined, instead no information is provided in these cases.

| Target compounds | CAS number | <i>tℝ</i> <sup>5</sup> , s | LOD <sup>2</sup> , ppm | LOQ⁴,ppm | Upper range³, ppm |
|------------------|------------|----------------------------|------------------------|----------|-------------------|
| Butanone, 2-     | 78-93-3    | 12.9                       | 1.00                   | 3.00     | 300               |
| Butyl acrylate   | 141-32-2   | 125.5                      | -                      | -        | -                 |
| Chlorobenzene    | 108-90-7   | 75.6                       | 1.00                   | 3.00     | 200               |
| Epichlorohydrin  | 106-89-8   | 27.3                       | 0.67                   | 2.00     | 200               |
| Ethyl acetate    | 141-78-6   | 14.6                       | 1.00                   | 3.00     | 300               |
| Ethyl acrylate   | 140-88-5   | 24.9                       | 1.00                   | 3.00     | 200               |

Qualified target compounds can be added to analysis programs and be assigned due to their retention time during analyses. The concentration calculation takes place using simplified assumptions as standards without the claim of high accuracy.

For more information see technical manual.

Technical manual Dräger X-pid® 8500/9000/9500 (draeger.com)

<sup>5</sup> Retention time