# PluraSens®



# Solvent Vapor Transmitter E2618-PID

**User Manual** 



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### Acetone

A colorless organic solvent with a characteristic pungent odor that is volatile and flammable with the chemical formula (CH<sub>3</sub>)<sub>2</sub>CO.

Synonyms/Trade Names: Dimethyl ketone, Ketone propane, 2-Propanone.

| Chemical formula                     |                            | (CH₃)₂CO  |  |  |
|--------------------------------------|----------------------------|---|--|--|
| Molar weight                         |                            | 58  |  |  |
| Relative gas den                     | sity (to air)              | 2.0   |  |  |
| Conversion*                          |                            | 1 ppm = 2.38 mg/m <sup>3</sup>  |  |  |
| Boiling point                        |                            | 56.11 °C  |  |  |
| Low explosive lir                    | mit (LEL), % vol. in air   | 2.5   |  |  |
| Upper explosive                      | limit (UEL), % vol. in air | 12.8  |  |  |
| Odor                                 |                            | Characteristic pungent smell  |  |  |
| Hazards                              |                            | Highly flammable.<br>Slightly toxic in normal use. Irritant causing mild skin<br>irritation and moderate to severe eye irritation. At high<br>vapor concentrations, it may depress the CNS. |  |  |
| Exposure limits 8 hours (2000/39/EC) |                            | 1900 mg/m³ / 500 ppm  |  |  |
|                                      | NIOSH REL TWA              | 590 mg/m³ /250 ppm  |  |  |
| IDLH (NIOSH)                         |                            | 2500 ppm [10%LEL]   |  |  |

### Benzene

A colorless liquid with an aromatic odor that is highly flammable with the molecular formula  $C_6H_6.$ 

Synonyms/Trade Names: Benzol, Phenyl hydride.

| Chemical formula   |                          | C <sub>6</sub> H <sub>6</sub>  |  |  |
|--------------------|--------------------------|--|--|--|
| Molar weight       |                          | 78   |  |  |
| Relative gas dens  | ity (to air)             | 2.69   |  |  |
| Conversion*        |                          | 1 ppm = 3.19 mg/m <sup>3</sup>   |  |  |
| Boiling point      |                          | 80 °C  |  |  |
| Low explosive lim  | it (LEL), % vol. in air  | 1.2  |  |  |
| Upper explosive li | mit (UEL), % vol. in air | 7.8  |  |  |
| Odor               |                          | Hyacinth-like odor   |  |  |
| Hazards            |                          | Highly flammable.<br>Irritant. Carcinogen.<br>May cause dizziness; headache, nausea, staggering<br>gait; anorexia, lassitude.<br>Target organs: eyes, skin, respiratory system, blood,<br>central nervous system, bone marrow. |  |  |
| Exposure limits    |                          | 0.319 mg/m³ /0.1 ppm   |  |  |
| (NIOSH REL)        | STEL<br>15 minutes       | 1 ppm  |  |  |
| Ca IDLH 500 pp     |                          |  |  |  |

# **Ethanol**

A clear colorless liquid with a faintly sweet odor and pungent taste.

Synonyms/Trade Names: ethyl, alcohol.

| Chemical formula                         |                     | CH₃CH₂OH  |
|--|---------------------|---|
| Molar weight                             |                     | 46  |
| Relative gas density (to                 | o air)              | 1,59  |
| Conversion*                              |                     | 1 ppm = 1.89 mg/m³  |
| Boiling point                            |                     | 78.37°C   |
| Low explosive limit (LE                  | EL), % vol. in air  | 3 - 3.3   |
| Upper explosive limit (                  | UEL), % vol. in air | 19  |
| Odor                                     |                     | The characteristic smell of alcohol   |
| Hazards                                  |                     | Highly flammable. Gas/air mixtures are explosive. Inhalation of vapors leads to cough, headache, fatigue, and drowsiness. High concentrations may damage the fetus. Repeated high exposure may affect the liver and the nervous system. |
| Exposure limits TWA according to 8 hours |                     | 1210 mg/m³ / 1000 ppm   |
| Commission Directive 2006/15/EC          | STEL<br>15 minutes  | -   |

# **Ethyl acetate**

A colorless organic compound with a sweet pear-like smell that is highly flammable.

Synonyms/Trade Names: ethyl ester, acetic ester, EA, EtOAc.

| Chemical formula                           |                | С <sub>4</sub> H <sub>8</sub> O <sub>2</sub>   |  |  |
|--|----------------|--|--|--|
| Molar weight                               |                | 88   |  |  |
| Conversion*                                |                | 1 ppm = 3.60 mg/m <sup>3</sup>   |  |  |
| Boiling point                              |                | 77.1 °C  |  |  |
| Low explosive limit (Lin air               | EL), % vol.    | 2  |  |  |
| Upper explosive limit (UEL), % vol. in air |                | 11.5   |  |  |
| Odor                                       |                | Sweet "pear" smell   |  |  |
| Hazards                                    |                | Flammable. Short-term exposure to high levels of ethyl acetate results first in irritation of the eyes, nose, and throat, followed by headache, nausea, vomiting, sleepiness, and unconsciousness. |  |  |
| Exposure limits (NIOSH)                    | TWA<br>8 hours | 1400 mg/m³ /400 ppm  |  |  |
|  | IDLH           | 2000 ppm [10%LEL]  |  |  |

# Toluene

A clear, colorless liquid with a sweet smell that is highly flammable.

Synonyms/Trade Names: methylbenzene, phenylmethane.

| Chemical formula                                      |                 | C <sub>6</sub> H₅CH₃ CH₃  |  |  |
|---|-----------------|---|--|--|
| Molar weight  |                 | 92  |  |  |
| Conversion*   |                 | 1 ppm = 3.77 mg/m <sup>3</sup>  |  |  |
| Boiling point   |                 | 110.7°C   |  |  |
| Low explosive limi in air                             | t (LEL), % vol. | 1.1 - 1.27  |  |  |
| Upper explosive limit (UEL), % vol. in air            |                 | 6.75-7.1  |  |  |
| Odor  |                 | Characteristic "chemical" smell   |  |  |
| Hazards   |                 | Highly flammable. Gas/air mixtures are explosive. Inhalation possible effects: irritation eyes, nose; lassitude (weakness, exhaustion), confusion, euphoria, dizziness, headache; dilated pupils, lacrimation (discharge of tears); anxiety, muscle fatigue, insomnia; paraesthesia; dermatitis; liver, kidney damage |  |  |
| Exposure limits                                       | TWA             | 192 mg/mm³ / 50 ppm   |  |  |
| according to<br>Commission<br>Directive<br>2006/15/EC | STEL            | 384 mg/mm³ / 100 ppm  |  |  |

# **Xylene**

A colorless, flammable liquid with a chemical smell. Xylenes are produced by the methylation of toluene and benzene.

The term is used for any of three isomers of dimethylbenzene, or a combination thereof.

Synonyms/Trade Names: xylol, dimethylbenzene.

| Chemical formula                         |                    |  |                           | C <sub>6</sub> H <sub>4</sub> (CH <sub>3</sub> ) <sub>2</sub> |  |
|--|--------------------|--|---------------------------|---|--|
| Isomers                                  |                    | ortho-xylene   | снь<br>снь<br>meta-xylene | ңс-Су-сң<br>para-xylene                                       |  |
| Molar weight                             |                    |  |                           | 106   |  |
| Conversion*                              |                    |  | 1 ppm = 4                 | 4.34 mg/m <sup>3</sup>  |  |
| Boiling point                            |                    | 144.4 °C   | 139 °C                    | 138.35 °C   |  |
| Low explosive limit (LEL), % vol. in air |                    | 0.9 - 1.1  |                           |   |  |
| Upper explosive lin                      | nit (UEL), % vol.  | 6.0-7.0  |                           |   |  |
| Odor                                     |                    | Charact  | eristic "cher             | nical" smell  |  |
| Hazards                                  |                    | Flammable. Inhaling can cause dizziness, headache, drowsiness, and nausea. |                           |   |  |
| Exposure limits according to             | TWA<br>8 hours     | 221 mg/mm³ / 50 ppm  |                           |   |  |
| Commission<br>Directive<br>2000/39/EC    | STEL<br>15 minutes | 442 mg/mm³ / 100 p   |                           |   |  |

Conversion of ppm to mg/m³ is calculated for 25°C and 1 atm.

# **Specifications**

| Detected gases            | V   | OCs with ionisation potential < 10.6 eV  |  |  |  |
|---------------------------|---|--|--|--|--|
| Sampling method           |   | Diffusion  |  |  |  |
| Sensor type               |   | Photoionization detector   |  |  |  |
| Calibration               |   | Isobutylene  |  |  |  |
| Order code                | E2618-PID-40  | E2618-PID-200  |  |  |  |
| Typical detection ranges  | 040 ppm (isobutylene)   | 0200 ppm (isobutylene)   |  |  |  |
| Resolution / digital unit | 1 ppb   | 1 ppm  |  |  |  |
| Response time T90         |   | <3s  |  |  |  |
| Sensor lifetime           |   | 5 years  |  |  |  |
| Calibration interval      | Monthly or more frequer   | ntly depending on operating conditions   |  |  |  |
| Signal update             |   | Every 1 second   |  |  |  |
| Load resistance           |   | $R_L$ < (Us - 2 V) / 22 mA for 4-20 mA $R_L$ > 250 kOhm for 0-10 V mode  |  |  |  |
| Digital interface         |   | RS485, Modbus RTU protocol<br>No galvanic isolation  |  |  |  |
| Power supply              | 1130 VDC or 24 VAC  |  |  |  |  |
| Power consumption         | < 2 VA  |  |  |  |  |
| Analog outputs            | 2 × 4-20 mA / 0-10 V, user settable   |  |  |  |  |
| Outputs assignment        |   | OUT1 2 gas; OUT2 2 gas   |  |  |  |
| Cable connections         |   | Screwless spring-loaded terminals  |  |  |  |
| Enclosure                 | Grey /  | ABS, wall mount, protection class IP65   |  |  |  |
| Dimensions                |   | H82 × W80 × D55 mm   |  |  |  |
| CE marking                | EN 61000-6-3:202<br>EN 61000-6-1:201  | J and 2014/35/EU, EN 50491-4-1:2012<br>20, EN 61326-1:2013(EMC, emissions)<br>9, EN 61000-6-2:2019(EMC, Immunity)<br>79-29-2:2015 and EN 60079-29-3:2014 |  |  |  |
| Operating conditions      | -40+65°C; 0,91,1 atm<br>095% RH non-condensing,<br>Explosion-safe indoor areas, Non-aggressive atmosphere |  |  |  |  |
| Other options             |   |  |  |  |  |
| Remote probe              | Protection IP65, default cable length 3.0 m;<br>max height 80 mm, max diameter 65 mm                      |  |  |  |  |

# **Product description**

E2618 series transmitters belong to the PluraSens® family of multifunctional measurement instruments. The instruments utilize gas sensors of various types with excellent repeatability, stability, and long lifetime.

E2618 series provide two independent analog outputs OUT1 and OUT2, user-selectable to 4-20 mA or 0-10 V, proportional to the chosen scale for gas concentration. RS485 Modbus RTU digital communication interface allows easy instrument configuration and integration into various automation systems.

Remote probe and duct mount versions are available.

The version of your transmitter is marked on the package. If the A symbol is marked on the product, consult the documentation for further information.

# Safety requirements

Misuse will impair the protection of the product. Always adhere to the safety provisions applicable in the country of use.

Do not perform any maintenance operation with the power on. Do not let water or foreign objects inside the device.

Removal of the PCB from the enclosure voids the warranty. Do not touch the electronic components directly, as they are sensitive to static electricity.

Connection diagrams can be found in the electrical connections section. The device might not perform correctly or be damaged if the wrong power supply is connected.

External circuits connected to the equipment should have sufficient insulation rating according to the environmental conditions and equipment power.

A disconnecting device that is marked as such and easily accessible should be included in the installation of this product.

# **Operating conditions**

The device should be used both in a non-hazardous area and in a basic electromagnetic environment, where the latter is defined in EN 61326-1. Avoid strong mechanical shock and vibrations. Avoid corrosive atmosphere and areas highly contaminated with dust, oil mist, etc. Keep the instrument away from direct sunlight. A sudden temperature or humidity change might affect the sensitivity of the sensor.

# Installation guidelines

Before proceeding with the installation it is mandatory to read the Safety requirements section and make sure to comply with all listed instructions. Installation standards EN 60079-29-2 and EN 45544-4 are also recommended for further instructions and related information about the installation. During the installation of the transmitter the following points must be considered:

- Application (air quality control or leakage detection.)
- Properties of the space under investigation (room geometry, direction, and velocity of airflows, etc.),
- Solvent vapors are heavier than air, so the gas detector-transmitter should be placed near the floor or potential leakage or formation source.
- The device should be accessible for maintenance and repair.

The aforementioned conditions above will affect the coverage area of the device. However, the coverage area for a transmitter is usually between 2.5 to 5 meters radius.

For early leakage detection, install the sensor as close as possible to the potential leakage sources (flanges, valves, pressure reducers, pumps, etc.), taking into consideration other points listed above.

For general area monitoring without definite leakage sources, the transmitters should be distributed evenly in the room.

For personal safety control, the transmitters are installed in the breathing zone (at the height of the head of people or animals). The recommended sensor position is vertical, pointing downwards.

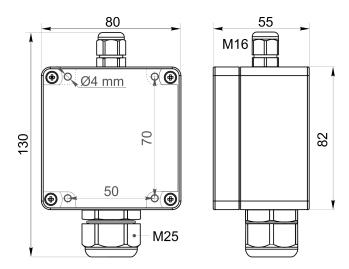
#### Wall mount version

Unscrew four lid screws and detach the lid from the transmitter. Attach the device to a wall with screws passing through mounting holes (for dimensions see the drawing in the next section).

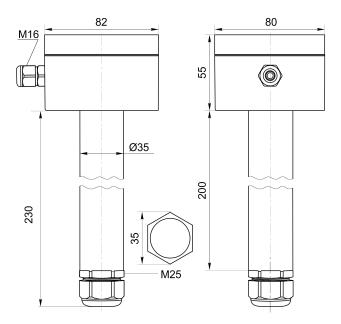
#### **Duct mount version**

Cut a hole with a diameter of 36...45 mm in the air duct at the chosen mounting place. Place the rubber flange aligning the holes in the flange and the air-duct and fix the flange with four self-tapping screws. Pass the sensor probe through the flange and adjust it to the appropriate depth. Unscrew four lid screws and detach the lid from the transmitter.

# **Mounting dimensions**



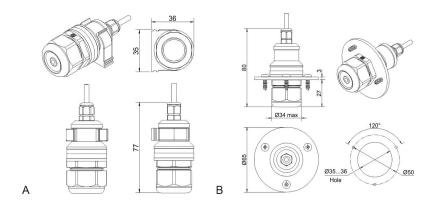
Wall mount version



Duct mount version

# Sensor probe handling

The wall-mount version of the detector-transmitter is available with a remote probe (see drawing below for dimensions). The remote probe is connected to the main unit with a shielded cable. The default remote probe cable length is 3 m.



- A) Wall mount remote probe with fixing clamp (default version),
- B) Remote probe with rubber flange and three self-tapping screws (on request)

The sensor probes of all types are equipped with a hydrophobic microporous PTFE filter to protect the sensor from dust, dirt, and water drops. The filter should be replaced if it gets strongly contaminated. To replace the PTFE filter, unscrew the sensor head cap and remove the old filter. Place a new filter into the cap and tighten it again.

**NOTE!** Never stab or press the filter near its center where the sensor is located since this may damage the sensor. Do not remove the filter as it may cause the device to show incorrect values and/or break the sensor.

The recommended orientation of the sensor probe is vertical with the sensor tip pointing downwards. This prevents the possible accumulation of condensed water on the sensor protection filter.

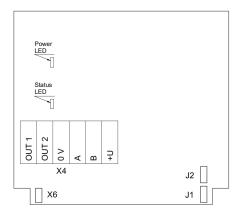
# Gas sensor replacement procedures

- 1. Remove the sensor head cap from the device (or the remote probe),
- 2. Remove the PTFE filter (if it is not removed within the cap),
- 3. Remove the O-ring rubber.
- 4. Detach the PID sensor from the device.
- 5. Insert the new PID sensor inside,
- 6. Put back the O-ring rubber, PTFE filter, and the head cap, respectively.

In most cases replacing the lamp and electrode stack can bring the sensor back to life.

### **Electrical connections**

Unscrew four lid screws and detach the lid from the device. Use the M16 cable gland to let in cables of the power supply and of the external devices. Attach the power cable to the device without turning it on. Using the connection diagram below, connect the analog outputs and digital interface terminals to the relevant devices according to your tasks.



PCB without PSU and relays

| Jumpers      |  |
|--------------|--|
| J1           | OUT1 type (open: 4-20 mA; closed 0-10 V)   |
| J2           | OUT2 type (open: 4-20 mA; closed 0-10 V)   |
| Х6           | Reset Modbus network parameters to default |
| X4 terminals |  |
| OUT1         | 4-20 mA / 0-10 V output                    |
| OUT2         | 4-20 mA / 0-10 V output                    |
| 0V           | 0 V / 24 VAC Neutral (optional)            |
| A            | RS485 A / Data +                           |
| В            | RS485 B / Data -                           |
| +U           | +24 VDC / 24 VAC Phase (optional)          |

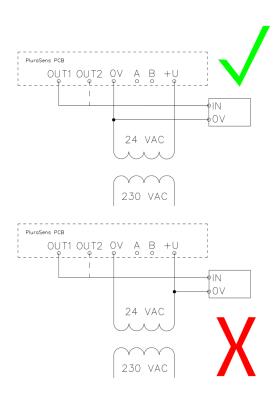
The screwless quick connect spring terminals on the E2618 series devices are suitable for a wide range of wires with a cross-section of 0.2...1.5 mm<sup>2</sup>. We recommend striping the wire end by 8...9 mm and using wire end sleeves.

To connect the wire, insert the wire end into the terminal hole. To disconnect, push the spring-loaded terminal lever, pull the wire out, and release the lever.

Use a twisted-pair cable, e.g. LiYY TP 2×2×0.5 mm<sup>2</sup> or CAT 5, to connect the device to the RS485 network. A and B on the device represent DATA+ and DATA- respectively, polarity must be respected when connecting to an external RS485 network.

Both analog outputs can be independently changed between 4-20 mA and 0-10 V type using jumpers J1 (OUT1) and J2 (OUT2). By closing pins on a specific jumper the related output is 0-10 V, with an open jumper the output is 4-20 mA. Power restart is required after changing the position of the jumpers.

#### Correct and incorrect cabling for 24 VAC



# Operation

Turn on the power. The instrument warm-up time takes ≤1 minute after switching on and the final sensor stabilization time to maximum accuracy takes <5 minutes. The operating status is indicated by the LED on the PCB of the device. The control LED (red) response to different processes is presented in the following table:

| Mode  | LED mode                 |
|---|--------------------------|
| During calibration mode or sensor heating period (if activated) | 0.5 Hz (50% on, 50% off) |
| During Modbus communication cycle                               | Short on-off pulses      |
| Normal operating/measurement                                    | Continuously on or off   |

Make sure that the transmitter is properly mounted, the external devices connected, power LED (green) on and control LED (red) constantly lit. Place the lid back and fix it with the screws. The device is ready to use.

# Properties of VOC: ionization energy, the correction factor

Sensors used in E2618-PID are calibrated using isobutylene, but the PID is a broadband VOC detector, with a sensitivity that differs for each VOC. If you know what VOC you are measuring, then the table below will allow you to calculate the concentration for your specific VOC. Only the most common substances are listed in the table, if you can't find your compound of concern, please contact us.

**NOTE!** These are approximate values, so for the best accuracy, you should calibrate with the relevant VOC.

The table includes the following columns:

- 1. The most common name for the VOC or other substances,
- 2. CAS No; You can find the VOC using the CAS No,
- Brutto-formula.
- 4. Ionization energy (IE), (eV)
- Relative Response/ Correction Factor (CF), Also called the Response Factor (RF).
   Multiply the displayed concentration by the Relative Response/CF/RF to calculate
   the actual concentration of the VOC.
- Minimum Detection Level (MDL) Also called Minimum Detectable Quantity (MDQ).
   The typical lowest concentration that can be detected. The sensor used in E2618-PID-40 has greater sensitivity than that in E2618-PID-200, so the MDL for the E2618-PID-40 will be much less than the MDL for the E2618-PID-200.

The Relative Response/CF/RF is measured in dry air; high humidity will reduce this factor by 30% to 50%, so the CF/RF should be increased in high humidities. Relative sensitivity is the inverse of the correction factor, specifying the percent response of the VOC, relative to isobutylene. If less than 100%, then the VOC is less responsive than isobutylene; if the relative sensitivity is greater than 100%, then the VOC is more responsive than isobutylene. Relative sensitivity (%) is specified in the same way as cross-sensitivity for toxic gas sensors.

#### **VOC** response

The PID can not measure all VOCs or gases, two types of VOCs are not measured:

NR: No response. The lamp does not ionize the VOC and the VOC cannot be measured.

NV: The vapor pressure of the VOC at 20°C is less than a few ppm, so this Semi-Volatile Organic Compound (SVOC) cannot be measured.

NA: Not available

Occasionally you will be measuring a mixture of VOCs. If the total concentration is within the linear range of your PID, then it is reasonable to assume that the concentrations are additive without interference between the different VOCs. Remember that if you are measuring a combination of VOCs, then the accurate measurement of one of these VOCs will be difficult; without careful data analysis, you will get only a CF averaged measurement. Be cautious when reporting actual VOC concentration if you know that there may be several VOCs present.

#### Balance gas

The relative response is measured in laboratory air, with 20.9% oxygen, balanced nitrogen. Some gases absorb UV light without causing any PID response (e.g. methane, ethane). In ambient atmospheres where these gases are present, the measured concentration of target gas will be less than is actually present.

Methane absorbs UV strongly, so for accurate measurements in methane containing atmospheres, calibrate with a calibration gas containing the expected methane concentration. 50% of LEL methane reduces the reading by up to 50%. Gases such as nitrogen and helium do not absorb UV and do not affect the relative response.

The correction factor for a gas mix containing PID detectable gases A, B, C... with response factors RF(A), RF(B), RF(C), in relative proportions a: b: c... is given by:

$$CF(mix) = 1 / [(a/CF(A) + b/CF(B) + c/CF(C)...]$$

Accuracy of the Table: This table is for indication only. Table accuracy is 1 to 2 digits only, so when calculating concentration for a specific VOC, only specify to 1 or 2 digits.

| Index | Chemical name    | Alternative name                                 | Formula                                      | CAS no.  | IE, eV | Response<br>Factor (RF)<br>10.6 eV | E2618-<br>PID-40<br>(ppb) | E2618-<br>PID-200<br>(ppb) |
|-------|------------------|--|--|----------|--------|------------------------------------|---------------------------|----------------------------|
| 1     | Acetaldehyde     |  | C <sub>2</sub> H <sub>4</sub> O              | 75-07-0  | 10.23  | 5.5                                | 25                        | 480                        |
| 2     | Acetamide        |  | C <sub>2</sub> H <sub>5</sub> NO             | 60-35-5  | 9.69   | 2                                  |                           |                            |
| 3     | Acetic acid      |  | C <sub>2</sub> H <sub>4</sub> O <sub>2</sub> | 64-19-7  | 10.66  | 28                                 | 180                       | 3615                       |
| 4     | Acetic anhydride |  | C <sub>4</sub> H <sub>6</sub> O <sub>3</sub> | 108-24-7 | 10.14  | 4                                  | 20                        | 400                        |
| 5     | Acetone          | 2-propanone,<br>dimethyl ketone                  | C <sub>3</sub> H <sub>6</sub> O              | 67-64-1  | 9.69   | 1.17                               | 5                         | 70                         |
| 6     | Acetonitrile     |  | CH₃CN  | 75-05-8  | 12.2   | NR                                 |                           |                            |
| 7     | Acetophenone     | 1-phenylethan-1<br>-one, methyl<br>phenyl ketone | C <sub>8</sub> H <sub>8</sub> O              | 98-86-2  | 9.29   | 0.6                                |                           |                            |
| 9     | Acetylene        | ethyne   | C <sub>2</sub> H <sub>2</sub>                | 74-86-2  | 11.4   | NR                                 |                           |                            |
| 10    | Acrolein         | propenal   | C₃H₄O  | 107-02-8 | 10.22  | 3.2                                | 20                        | 400                        |

| 11    | Acrylic Acid             | propenoic acid                  | C <sub>3</sub> H <sub>4</sub> O <sub>2</sub>   | 79-10-7   | 10.6   | 21                                 | 15                        | 275                        |
|-------|--------------------------|---------------------------------|--|-----------|--------|------------------------------------|---------------------------|----------------------------|
| Index | Chemical name            | Alternative name                | Formula  | CAS no.   | IE, eV | Response<br>Factor (RF)<br>10.6 eV | E2618-<br>PID-40<br>(ppb) | E2618-<br>PID-200<br>(ppb) |
| 12    | Acrylonitrile            |                                 | C <sub>3</sub> H <sub>3</sub> N                | 107-13-1  | 10.91  | NR                                 |                           |                            |
| 13    | Alkanes, n-, C6+         |                                 | C <sub>n</sub> H <sub>2n+2</sub>               | N/A       | ~10    | 1.2                                |                           |                            |
| 15    | Allyl alcohol            |                                 | C <sub>3</sub> H <sub>6</sub> O                | 107-18-6  | 9.63   | 2.3                                | 10                        | 200                        |
| 16    | Allyl bromide            | 3-bromopropene                  | C₃H₅Br   | 106-95-6  | 9.96   | 3                                  |                           |                            |
| 17    | Allyl chloride           | 3-chloropropene                 | C₃H₅CI   | 107-05-1  | 10.05  | 4.5                                | 20                        | 450                        |
| 18    | Ammonia                  |                                 | NH <sub>3</sub>                                | 7664-41-7 | 10.18  | 8.5                                | 40                        | 850                        |
| 19    | Amyl acetate             |                                 | C <sub>7</sub> H <sub>14</sub> O <sub>2</sub>  | 628-63-7  | 9.9    | 1.8                                | 10                        | 180                        |
| 20    | Amyl alcohol             |                                 | C <sub>5</sub> H <sub>12</sub> O               | 71-41-0   | 10     | 2.6                                | 15                        | 320                        |
| 21    | Amyl alcohol,tert-       |                                 | C <sub>5</sub> H <sub>12</sub> O               | 75-85-4   | 9.8    | 1.5                                |                           |                            |
| 22    | Aniline                  |                                 | C <sub>6</sub> H <sub>7</sub> N                | 62-53-3   | 7.7    | 0.5                                | 3                         | 50                         |
| 23    | Anisole                  |                                 | C <sub>7</sub> H <sub>8</sub> O                | 100-66-3  | 8.21   | 0.59                               | 2                         | 50                         |
| 24    | Arsine                   |                                 | AsH <sub>3</sub>                               | 7784-42-1 | 9.89   | 2.5                                | 15                        | 250                        |
| 25    | Asphalt, petroleum fumes |                                 |  | 8052-42-4 | ~9     | 1                                  | 5                         | 100                        |
| 26    | Benzaldehyde             |                                 | C <sub>7</sub> H <sub>6</sub> O                | 100-52-7  | 9.49   | 0.7                                | 5                         | 85                         |
| 27    | Benzene                  |                                 | C <sub>6</sub> H <sub>6</sub>                  | 71-43-2   | 9.24   | 0.5                                | 3                         | 50                         |
| 28    | Benzoic acid             |                                 | C <sub>7</sub> H <sub>6</sub> O <sub>2</sub>   | 65-85-0   | 9.3    | 0.7                                |                           |                            |
| 29    | Benzyl acetate           |                                 | C <sub>9</sub> H <sub>10</sub> O <sub>2</sub>  | 140-11-4  | ~9     | 0.6                                |                           |                            |
| 30    | Benzyl alcohol           |                                 | C <sub>7</sub> H <sub>8</sub> O                | 100-51-6  | 8.26   | 1                                  | 6                         | 125                        |
| 31    | Camphor                  |                                 | C <sub>10</sub> H <sub>16</sub> O              | 76-22-2   | 8.76   | 0.4                                |                           |                            |
| 32    | Carbon dioxide           |                                 | CO <sub>2</sub>                                | 124-38-9  | 13.77  | NR                                 | 0                         | 0                          |
| 33    | Carbon disulfide         |                                 | CS <sub>2</sub>                                | 75-15-0   | 10.08  | 1.4                                | 7                         | 140                        |
| 34    | Carbon monoxide          |                                 | со   | 630-08-0  | 14.01  | NR                                 | 0                         | 0                          |
| 35    | Carbon<br>tetrabromide   | tetrabromometh<br>ane           | CBr <sub>4</sub>                               | 558-13-4  | 10.31  | 3                                  | 15                        | 300                        |
| 36    | Carbon<br>tetrachloride  | R-10,<br>tetrachlorometha<br>ne | CCI <sub>4</sub>                               | 56-23-5   | 11.47  | NR                                 |                           |                            |
| 37    | Chloroform               |                                 | CHCl₃  | 67-66-3   | 11.42  | NR                                 |                           |                            |
| 38    | Acetaldehyde             |                                 | C <sub>14</sub> H <sub>26</sub> O <sub>2</sub> | 97-89-2   | ~9     | 0.9                                |                           |                            |
| 39    | Coumarin                 |                                 | C <sub>9</sub> H <sub>6</sub> O <sub>2</sub>   | 91-64-5   | ~9     | 0.4                                |                           |                            |
| 40    | Creosote                 |                                 | n/a  | 8021-39-4 | ~9     | 1                                  |                           |                            |
| 41    | Cresol, m-               | 3-methylphenol                  | C <sub>7</sub> H <sub>8</sub> O                | 108-39-4  | 8.36   | 2.2                                | 5                         | 105                        |

| 42    | Cresol, o-       | 2-methylphenol           | C <sub>7</sub> H <sub>8</sub> O               | 95-48-7    | 8.14   | 1.1                                | 5                         | 105                        |
|-------|------------------|--------------------------|---|------------|--------|------------------------------------|---------------------------|----------------------------|
| 43    | Cresol, p-       | 4-methylphenol           | C <sub>7</sub> H <sub>8</sub> O               | 106-44-5   | 8.31   | 1.1                                | 5                         | 105                        |
| 44    | Cycloalkanes     |                          | N/A   | N/A        | ~10    | 1.5                                |                           |                            |
| 45    | Cyclobutene      |                          | C₄H <sub>6</sub>                              | 822-35-5   | 9.43   | 3                                  |                           |                            |
| 46    | Cycloheptane     |                          | C <sub>7</sub> H <sub>14</sub>                | 291-64-5   | 9.82   | 1.1                                |                           |                            |
| 47    | Cyclohexane      |                          | C <sub>6</sub> H <sub>12</sub>                | 110-82-7   | 9.98   | 1.3                                | 7                         | 130                        |
| 48    | Cyclohexanethiol |                          | C <sub>6</sub> H <sub>12</sub> S              | 1569-69-3  | ~9     | 0.5                                |                           |                            |
| 49    | Cyclohexanol     |                          | C <sub>6</sub> H <sub>12</sub> O              | 108-93-0   | 10     | 1.6                                | 15                        | 300                        |
| 50    | Cyclohexene      |                          | C <sub>6</sub> H <sub>10</sub>                | 110-83-8   | 8.95   | 0.9                                | 5                         | 75                         |
| 51    | Cyclopentadiene  |                          | C₅H <sub>6</sub>                              | 542-92-7   | 8.56   | 0.8                                |                           |                            |
| Index | Chemical name    | Alternative name         | Formula                                       | CAS no.    | IE, eV | Response<br>Factor (RF)<br>10.6 eV | E2618-<br>PID-40<br>(ppb) | E2618-<br>PID-200<br>(ppb) |
| 52    | Cyclopentane     |                          | C₅H₁0   | 287-92-3   | 10.52  | 10                                 | 20                        | 400                        |
| 53    | Cyclopentene     |                          | C₅H <sub>8</sub>                              | 142-29-0   | 9.01   | 1.5                                |                           |                            |
| 54    | Decane, n-       |                          | C <sub>10</sub> H <sub>22</sub>               | 124-18-5   | 9.65   | 1.2                                | 5                         | 100                        |
| 55    | Decanol          |                          | C <sub>10</sub> H <sub>22</sub> O             | 112-30-1   |        | 1.2                                |                           |                            |
| 56    | Diethyl ether    | ethyl ether              | C <sub>4</sub> H <sub>10</sub> O              | 60-29-7    | 9.53   | 1.1                                | 4                         | 90                         |
| 57    | Diethylamine     |                          | C <sub>4</sub> H <sub>11</sub> N              | 109-89-7   | 8.01   | 1.4                                | 5                         | 100                        |
| 58    | Diesel fuel      |                          |   | 68334-30-5 | 8      | 0.8                                | 4                         | 75                         |
| 59    | Dimethylamine    |                          | C <sub>2</sub> H <sub>7</sub> N               | 124-40-3   | 8.24   | 1.5                                | 7                         | 140                        |
| 60    | Diphenyl ether   | phenyl ether             | C <sub>12</sub> H <sub>10</sub> O             | 101-84-8   | 8.09   | 1.5                                | 4                         | 80                         |
| 61    | Dodecane         |                          | C <sub>12</sub> H <sub>24</sub>               | 112-40-3   | ~8.8   | 1                                  |                           |                            |
| 62    | Dodecanol        |                          | C <sub>12</sub> H <sub>26</sub> O             | 112-53-8   |        | 0.9                                | 5                         | 90                         |
| 63    | Ethane           |                          | C <sub>2</sub> H <sub>6</sub>                 | 74-84-0    | 11.56  | NR                                 |                           |                            |
| 64    | Ethanol          | alcohol,ethyl<br>alcohol | C <sub>2</sub> H <sub>6</sub> O               | 64-17-5    | 10.43  | 11                                 | 45                        | 870                        |
| 65    | Ethyl acetate    |                          | C <sub>4</sub> H <sub>8</sub> O <sub>2</sub>  | 141-78-6   | 10.01  | 4.5                                | 20                        | 360                        |
| 66    | Ethyl acrylate   |                          | C <sub>5</sub> H <sub>8</sub> O <sub>2</sub>  | 140-88-5   | 10.3   | 2.3                                | 10                        | 200                        |
| 67    | Ethyl benzoate   |                          | C <sub>9</sub> H <sub>10</sub> O <sub>2</sub> | 93-89-0    | 8.9    | 0.9                                |                           |                            |
| 68    | Ethyl butyrate   |                          | C <sub>6</sub> H <sub>12</sub> O <sub>2</sub> | 105-54-4   | ~9.9   | 1.4                                | 5                         | 100                        |
| 69    | Ethylamine       |                          | C <sub>2</sub> H <sub>7</sub> N               | 75-04-7    | 8.86   | 1                                  | 5                         | 100                        |
| 70    | Ethylbenzene     |                          | C <sub>8</sub> H <sub>10</sub>                | 100-41-4   | 8.76   | 0.56                               | 3                         | 50                         |
| 71    | Ethylcyclohexane |                          | C <sub>8</sub> H <sub>16</sub>                | 1678-91-7  | 9.54   | 0.8                                |                           |                            |
| 72    | Ethylene         | ethene                   | C <sub>2</sub> H <sub>4</sub>                 | 74-85-1    | 10.51  | 8                                  | 40                        | 800                        |
| 73    | Ethylene glycol  |                          | C <sub>2</sub> H <sub>6</sub> O <sub>2</sub>  | 107-21-1   | 10.16  | 9                                  | 100                       | 2000                       |
| 74    | Ethylamine       |                          | C <sub>2</sub> H <sub>7</sub> N               | 75-04-7    | 8.86   | 1                                  | 5                         | 100                        |

| 75  | Formaldehyde   |                   | CH₂O   | 50-00-0  | 10.87   | NR  |                                    |  |
|---|--|-------------------|--|--|---|---|------------------------------------|--|
| 76  | Formamide  |                   | CH₃ON  | 75-12-7  | 10.2  | 2   | 10                                 | 200  |
| 77  | Formic acid  |                   | CH <sub>2</sub> O <sub>2</sub>   | 64-18-6  | 11.05   | NR  |                                    |  |
| 78  | Furan  |                   | C <sub>4</sub> H <sub>4</sub> O  | 110-00-9   | 8.88  | 0.4   |                                    |  |
| 79  | Hexanoic acid  |                   | C <sub>6</sub> H <sub>12</sub> O <sub>2</sub>  | 142-62-1   | 10.12   | 4   |                                    |  |
| 80  | Hexanol  |                   | C <sub>6</sub> H <sub>14</sub> O   | 111-27-3   | 9.89  | 2   |                                    |  |
| 81  | Hexene, 1-   |                   | C <sub>6</sub> H <sub>12</sub>   | 592-41-6   | 9.44  | 0.98  | 5                                  | 90   |
| 82  | Hydrazine  |                   | H <sub>4</sub> N <sub>2</sub>  | 302-01-2   | 8.93  | 3   | 15                                 | 300  |
| 83  | Hydrazoic acid   |                   | HN <sub>3</sub>  | 7782-79-8  | 10.72   | NR  |                                    |  |
| 84  | Hydrogen   |                   | H <sub>2</sub>   | 1333-74-0  | 15.43   | NR  |                                    |  |
| 85  | Hydrogen bromide   |                   | HBr  | 10035-10-6   | 11.62   | NR  |                                    |  |
| 86  | Hydrogen chloride  |                   | HCI  | 7647-01-0  | 12.74   | NR  |                                    |  |
| 87  | Hydrogen cyanide   |                   | HCN  | 74-90-8  | 13.6  | NR  |                                    |  |
| 88  | Hydrogen fluoride  |                   | HF   | 7664-39-3  | 15.98   | NR  |                                    |  |
| 89  | Hydrogen iodide  |                   | НІ   | 10034-85-2   | 10.39   | 5   |                                    |  |
| 90  | Hydrogen peroxide  |                   | H <sub>2</sub> O <sub>2</sub>  | 7722-84-1  | 10.58   | 4   | 20                                 | 400  |
| Index   | Chemical name  | Alternative name  | Formula  | CAS no.  | IE, eV  | Response<br>Factor (RF)<br>10.6 eV                              | E2618-<br>PID-40<br>(ppb)          | E2618-<br>PID-200<br>(ppb)                   |
| 106   |  |                   |  |  |   |   |                                    |  |
| 126   | Hydrogen selenide  |                   | H₂Se   | 7783-07-5  | 9.88  | 2   |                                    |  |
| 126   | Hydrogen selenide Hydrogen sulfide   |                   | H <sub>2</sub> Se<br>H <sub>2</sub> S  | 7783-07-5<br>7783-06-4   | 9.88  | 2   |                                    |  |
|   | , ,  |                   |  |  |   |   | 4                                  | 80   |
| 127   | Hydrogen sulfide   | 2,2,4-trimethylpe | H₂S  | 7783-06-4  | 10.46   | 4   | 4 5                                | 80   |
| 127<br>128  | Hydrogen sulfide Hydroquinone  | ' '               | H <sub>2</sub> S<br>C <sub>6</sub> H <sub>6</sub> O <sub>2</sub>   | 7783-06-4<br>123-31-9  | 10.46<br>7.94<br>9.86   | 4<br>0.8  |                                    |  |
| 127<br>128<br>129   | Hydrogen sulfide Hydroquinone Isooctane  | ' '               | H <sub>2</sub> S<br>C <sub>6</sub> H <sub>6</sub> O <sub>2</sub><br>C <sub>8</sub> H <sub>18</sub>   | 7783-06-4<br>123-31-9<br>540-84-1  | 10.46<br>7.94<br>9.86   | 4<br>0.8<br>1.1   | 5                                  | 100  |
| 127<br>128<br>129<br>130  | Hydrogen sulfide Hydroquinone Isooctane Isooctanol   | ' '               | H <sub>2</sub> S<br>C <sub>6</sub> H <sub>6</sub> O <sub>2</sub><br>C <sub>8</sub> H <sub>18</sub><br>C <sub>8</sub> H <sub>18</sub> O   | 7783-06-4<br>123-31-9<br>540-84-1<br>26952-21-6  | 10.46<br>7.94<br>9.86<br>~9.8   | 4<br>0.8<br>1.1<br>1.7  | 5<br>9                             | 100<br>170                                   |
| 127<br>128<br>129<br>130<br>131   | Hydrogen sulfide Hydroquinone Isooctane Isooctanol Isopentane  | ' '               | H <sub>2</sub> S<br>C <sub>6</sub> H <sub>6</sub> O <sub>2</sub><br>C <sub>8</sub> H <sub>18</sub><br>C <sub>8</sub> H <sub>18</sub> O<br>C <sub>5</sub> H <sub>12</sub><br>C <sub>5</sub> H <sub>12</sub> O<br>C <sub>5</sub> H <sub>10</sub>   | 7783-06-4<br>123-31-9<br>540-84-1<br>26952-21-6<br>78-78-4   | 10.46<br>7.94<br>9.86<br>~9.8<br>10.32<br>9.86<br>9.12                                  | 4<br>0.8<br>1.1<br>1.7<br>4                                     | 5<br>9                             | 100<br>170                                   |
| 127<br>128<br>129<br>130<br>131<br>132                                    | Hydrogen sulfide Hydroquinone Isooctane Isooctanol Isopentane Isopentanol  | ' '               | H <sub>2</sub> S<br>C <sub>6</sub> H <sub>6</sub> O <sub>2</sub><br>C <sub>8</sub> H <sub>18</sub><br>C <sub>8</sub> H <sub>18</sub> O<br>C <sub>5</sub> H <sub>12</sub><br>C <sub>5</sub> H <sub>12</sub> O   | 7783-06-4<br>123-31-9<br>540-84-1<br>26952-21-6<br>78-78-4<br>137-32-6   | 10.46<br>7.94<br>9.86<br>~9.8<br>10.32<br>9.86  | 4<br>0.8<br>1.1<br>1.7<br>4<br>2                                | 5<br>9                             | 100<br>170                                   |
| 127<br>128<br>129<br>130<br>131<br>132<br>133                             | Hydrogen sulfide Hydroquinone Isooctane Isooctanol Isopentane Isopentanol Isopentene Iodoform Iodomethane  | ntane             | H <sub>2</sub> S<br>C <sub>6</sub> H <sub>6</sub> O <sub>2</sub><br>C <sub>8</sub> H <sub>18</sub><br>C <sub>8</sub> H <sub>18</sub> O<br>C <sub>5</sub> H <sub>12</sub><br>C <sub>5</sub> H <sub>12</sub> O<br>C <sub>5</sub> H <sub>10</sub>   | 7783-06-4<br>123-31-9<br>540-84-1<br>26952-21-6<br>78-78-4<br>137-32-6<br>563-46-2   | 10.46<br>7.94<br>9.86<br>~9.8<br>10.32<br>9.86<br>9.12                                  | 4<br>0.8<br>1.1<br>1.7<br>4<br>2<br>0.8                         | 5<br>9<br>30                       | 100<br>170<br>600                            |
| 127<br>128<br>129<br>130<br>131<br>132<br>133<br>134                      | Hydrogen sulfide Hydroquinone Isooctane Isooctanol Isopentane Isopentanol Isopentene Iodoform  | ntane             | H <sub>2</sub> S<br>C <sub>6</sub> H <sub>6</sub> O <sub>2</sub><br>C <sub>8</sub> H <sub>18</sub><br>C <sub>8</sub> H <sub>18</sub> O<br>C <sub>5</sub> H <sub>12</sub><br>C <sub>5</sub> H <sub>12</sub> O<br>C <sub>5</sub> H <sub>10</sub><br>CHI <sub>3</sub>   | 7783-06-4<br>123-31-9<br>540-84-1<br>26952-21-6<br>78-78-4<br>137-32-6<br>563-46-2<br>75-47-8  | 10.46<br>7.94<br>9.86<br>~9.8<br>10.32<br>9.86<br>9.12<br>9.25                          | 4<br>0.8<br>1.1<br>1.7<br>4<br>2<br>0.8<br>1.5                  | 5<br>9<br>30                       | 100<br>170<br>600                            |
| 127<br>128<br>129<br>130<br>131<br>132<br>133<br>134<br>135               | Hydrogen sulfide Hydroquinone Isooctane Isooctanol Isopentane Isopentanol Isopentene Iodoform Iodomethane Isoalkanes,                              | ntane             | H <sub>2</sub> S<br>C <sub>6</sub> H <sub>6</sub> O <sub>2</sub><br>C <sub>8</sub> H <sub>18</sub><br>C <sub>8</sub> H <sub>18</sub> O<br>C <sub>5</sub> H <sub>12</sub><br>C <sub>5</sub> H <sub>12</sub> O<br>C <sub>5</sub> H <sub>10</sub><br>CHI <sub>3</sub><br>CH <sub>3</sub> I<br>C <sub>8</sub> H <sub>18</sub> O<br>C <sub>4</sub> H <sub>10</sub>  | 7783-06-4<br>123-31-9<br>540-84-1<br>26952-21-6<br>78-78-4<br>137-32-6<br>563-46-2<br>75-47-8<br>74-88-4                                     | 10.46<br>7.94<br>9.86<br>~9.8<br>10.32<br>9.86<br>9.12<br>9.25<br>9.54                  | 4<br>0.8<br>1.1<br>1.7<br>4<br>2<br>0.8<br>1.5<br>0.4           | 5<br>9<br>30                       | 100<br>170<br>600                            |
| 127<br>128<br>129<br>130<br>131<br>132<br>133<br>134<br>135               | Hydrogen sulfide Hydroquinone Isooctane Isooctanol Isopentane Isopentanol Isopentene Iodoform Iodomethane Isoalkanes, C10-C13                      | ntane             | H <sub>2</sub> S<br>C <sub>6</sub> H <sub>6</sub> O <sub>2</sub><br>C <sub>8</sub> H <sub>18</sub><br>C <sub>8</sub> H <sub>18</sub> O<br>C <sub>5</sub> H <sub>12</sub><br>C <sub>5</sub> H <sub>12</sub> O<br>C <sub>5</sub> H <sub>10</sub><br>CHI <sub>3</sub><br>CH <sub>3</sub> I<br>C <sub>8</sub> H <sub>18</sub> O  | 7783-06-4<br>123-31-9<br>540-84-1<br>26952-21-6<br>78-78-4<br>137-32-6<br>563-46-2<br>75-47-8<br>74-88-4<br>68551-17-7                       | 10.46<br>7.94<br>9.86<br>~9.8<br>10.32<br>9.86<br>9.12<br>9.25<br>9.54<br>~9.6          | 4<br>0.8<br>1.1<br>1.7<br>4<br>2<br>0.8<br>1.5<br>0.4           | 5<br>9<br>30<br>8<br>2             | 100<br>170<br>600<br>150<br>40               |
| 127<br>128<br>129<br>130<br>131<br>132<br>133<br>134<br>135<br>136        | Hydrogen sulfide Hydroquinone Isooctane Isooctanol Isopentane Isopentanol Isopentene Iodoform Iodomethane Isoalkanes, C10-C13 Isobutane            | ntane             | H <sub>2</sub> S<br>C <sub>6</sub> H <sub>6</sub> O <sub>2</sub><br>C <sub>8</sub> H <sub>18</sub><br>C <sub>8</sub> H <sub>18</sub> O<br>C <sub>5</sub> H <sub>12</sub><br>C <sub>5</sub> H <sub>12</sub> O<br>C <sub>5</sub> H <sub>10</sub><br>CHI <sub>3</sub><br>CH <sub>3</sub> I<br>C <sub>8</sub> H <sub>18</sub> O<br>C <sub>4</sub> H <sub>10</sub><br>C <sub>4</sub> H <sub>10</sub> O<br>C <sub>6</sub> H <sub>12</sub> O <sub>2</sub> | 7783-06-4<br>123-31-9<br>540-84-1<br>26952-21-6<br>78-78-4<br>137-32-6<br>563-46-2<br>75-47-8<br>74-88-4<br>68551-17-7<br>75-28-5            | 10.46<br>7.94<br>9.86<br>~9.8<br>10.32<br>9.86<br>9.12<br>9.25<br>9.54<br>~9.6<br>10.57 | 4<br>0.8<br>1.1<br>1.7<br>4<br>2<br>0.8<br>1.5<br>0.4<br>1      | 5<br>9<br>30<br>8<br>2             | 100<br>170<br>600<br>150<br>40               |
| 127<br>128<br>129<br>130<br>131<br>132<br>133<br>134<br>135<br>136<br>137 | Hydrogen sulfide Hydroquinone Isooctane Isooctanol Isopentane Isopentanol Isopentene Iodoform Iodomethane Isoalkanes, C10-C13 Isobutane Isobutanol | ntane             | H <sub>2</sub> S<br>C <sub>6</sub> H <sub>6</sub> O <sub>2</sub><br>C <sub>8</sub> H <sub>18</sub><br>C <sub>8</sub> H <sub>18</sub> O<br>C <sub>5</sub> H <sub>12</sub><br>C <sub>5</sub> H <sub>12</sub> O<br>CHI <sub>3</sub><br>CH <sub>3</sub> I<br>C <sub>8</sub> H <sub>18</sub> O<br>C <sub>4</sub> H <sub>10</sub><br>C <sub>4</sub> H <sub>10</sub> O  | 7783-06-4<br>123-31-9<br>540-84-1<br>26952-21-6<br>78-78-4<br>137-32-6<br>563-46-2<br>75-47-8<br>74-88-4<br>68551-17-7<br>75-28-5<br>78-83-1 | 10.46<br>7.94<br>9.86<br>~9.8<br>10.32<br>9.86<br>9.12<br>9.25<br>9.54<br>~9.6<br>10.57 | 4<br>0.8<br>1.1<br>1.7<br>4<br>2<br>0.8<br>1.5<br>0.4<br>1<br>8 | 5<br>9<br>30<br>8<br>2<br>40<br>20 | 100<br>170<br>600<br>150<br>40<br>800<br>350 |

| 142   | Isooctanol                    |  | C <sub>8</sub> H <sub>18</sub> O              | 26952-21-6 | ~9.8      | 1.7                                | 9                         | 170                        |
|-------|-------------------------------|--|---|------------|-----------|------------------------------------|---------------------------|----------------------------|
| 143   | Isopentane                    |  | C <sub>5</sub> H <sub>12</sub>                | 78-78-4    | 10.32     | 4                                  | 30                        | 600                        |
| 144   | Isopentanol                   |  | C <sub>5</sub> H <sub>12</sub> O              | 137-32-6   | 9.86      | 2                                  |                           |                            |
| 145   | Isopentene                    |  | C <sub>5</sub> H <sub>10</sub>                | 563-46-2   | 9.12      | 0.8                                |                           |                            |
| 146   | Isoprene                      | 2-methyl-1,3-buta diene                | C₅H <sub>8</sub>                              | 78-79-5    | 8.85      | 0.8                                | 3                         | 70                         |
| 147   | Isopropanol                   | IPA, 2-propanol                        | C₃H <sub>8</sub> O                            | 67-63-0    | 10.17     | 4                                  | 22                        | 440                        |
| 148   | Isovaleraldehyde              |  | C <sub>5</sub> H <sub>10</sub> O              | 590-86-3   | 9.72      | 1.3                                |                           |                            |
| 149   | Isovaleric Acid               |  | C <sub>5</sub> H <sub>10</sub> O <sub>2</sub> | 503-74-2   | ~10.<br>2 | 5.5                                |                           |                            |
| 150   | Methane                       | natural gas                            | CH₄   | 74-82-8    | 12.51     | NR                                 |                           |                            |
| 151   | Methanol                      |  | CH₄O  | 67-56-1    | 10.85     | NR                                 | 1000                      | 20000                      |
| 152   | Methyl ethyl ketone           | MEK,<br>Butan-2-one                    | C₄H <sub>8</sub> O                            | 78-93-3    | 9.51      | 0.96                               | 4                         | 80                         |
| 153   | Methyl ethyl ketone peroxides | MEKP                                   | C <sub>8</sub> H <sub>18</sub> O <sub>6</sub> | 1338-23-4  | ~9        | 0.8                                | 4                         | 80                         |
| 154   | Methyl formate                |  | C <sub>2</sub> H <sub>4</sub> O <sub>2</sub>  | 107-31-3   | 10.82     | NR                                 |                           |                            |
| 155   | Methyl mercaptan              |  | CH₄S  | 74-93-1    | 9.44      | 0.7                                | 4                         | 70                         |
| 156   | Methylamine                   |  | CH₅N  | 74-89-5    | 8.97      | 1.5                                | 7                         | 140                        |
| 157   | Mineral oil                   |  |   | 8042-47-5  | ~9        | 0.8                                | 4                         | 80                         |
| 158   | Mineral spirits               | Stoddard<br>solvent, Varsol,<br>Viscor |   | 64475-85-0 | ~9        | 0.8                                | 4                         | 80                         |
| 159   | Naphthalene                   |  | C <sub>10</sub> H <sub>8</sub>                | 91-20-3    | 8.14      | 0.4                                | 2                         | 45                         |
| 160   | Neopentane                    | tetramethylmeth<br>ane                 | C <sub>5</sub> H <sub>12</sub>                | 207-343-7  | 10.21     | 3                                  |                           |                            |
| 161   | Neopentyl alcohol             |  | C <sub>5</sub> H <sub>12</sub> O              | 75-84-3    | 9.72      | 2                                  |                           |                            |
| 162   | Nitric oxide                  |  | NO  | 10102-43-9 | 9.27      | 8                                  | 40                        | 800                        |
| 163   | Nitrobenzene                  |  | C <sub>6</sub> H <sub>5</sub> NO <sub>2</sub> | 98-95-3    | 9.92      | 1.7                                | 10                        | 170                        |
| 164   | Nitroethane                   |  | C <sub>2</sub> H <sub>5</sub> NO <sub>2</sub> | 79-24-3    | 10.88     | NR                                 |                           |                            |
| Index | Chemical name                 | Alternative name                       | Formula                                       | CAS no.    | IE, eV    | Response<br>Factor (RF)<br>10.6 eV | E2618-<br>PID-40<br>(ppb) | E2618-<br>PID-200<br>(ppb) |
| 165   | Nitrogen trichloride          |  | NCl <sub>3</sub>                              | 10025-85-1 | 10.1      | 1                                  | 5                         | 100                        |
| 166   | Nitrogen                      |  | $N_2$   | 7727-37-9  | 15.58     | NR                                 |                           |                            |
| 167   | Nitrogen dioxide              |  | NO <sub>2</sub>                               | 10102-44-0 | 9.58      | 10                                 | 50                        | 1000                       |
| 168   | Nitrogen trifluoride          |  | NF <sub>3</sub>                               | 7783-54-2  | 12.97     | NR                                 |                           |                            |
| 169   | Nitromethane                  |  | CH <sub>3</sub> NO <sub>2</sub>               | 75-52-5    | 11.08     | NR                                 |                           |                            |
| 170   | Nitropropane, 1-              |  | C <sub>3</sub> H <sub>7</sub> NO <sub>2</sub> | 108-03-2   | 10.81     | NR                                 |                           |                            |
| 171   | Nitropropane, 2-              |  | C <sub>3</sub> H <sub>7</sub> NO <sub>2</sub> | 79-46-9    | 10.71     | NR                                 |                           |                            |

|     | 1                       |                                   |   |            |           |      |    |      |
|-----|-------------------------|-----------------------------------|---|------------|-----------|------|----|------|
| 172 | Nitrous oxide           |                                   | N <sub>2</sub> O                              | 10024-97-2 | 12.89     | NR   |    |      |
| 173 | Nonane                  |                                   | C <sub>9</sub> H <sub>20</sub>                | 111-84-2   | 9.72      | 1.4  | 6  | 130  |
| 174 | Nonanol (mixed isomers) |                                   | C <sub>9</sub> H <sub>20</sub> O              | 143-08-8   | ~9.8      | 1.2  |    |      |
| 175 | Nonene (mixed isomers)  |                                   | C <sub>9</sub> H <sub>18</sub>                | 27215-95-8 | ~9.3      | 0.6  |    |      |
| 176 | Nonene, 1-              |                                   | C <sub>9</sub> H <sub>18</sub>                | 124-11-8   | ~9.4      | 0.6  |    |      |
| 177 | Octane                  |                                   | C <sub>8</sub> H <sub>18</sub>                | 111-65-9   | 9.8       | 1.6  | 8  | 160  |
| 178 | Octanol (mixed isomers) | capryl alcohol,<br>octyl alcohol  | C <sub>8</sub> H <sub>18</sub> O              | 111-87-5   | ~9.8      | 1.5  |    |      |
| 179 | Octene (mixed isomers)  |                                   | C <sub>8</sub> H <sub>16</sub>                | 25377-83-7 | ~9.4      | 0.7  |    |      |
| 180 | Octene, 1-              |                                   | C <sub>8</sub> H <sub>16</sub>                | 111-66-0   | 9.43      | 0.7  | 3  | 70   |
| 181 | Oxalic acid             |                                   | C <sub>2</sub> H <sub>2</sub> O <sub>4</sub>  | 144-62-7   | 11.2      | NR   |    |      |
| 182 | Oxalyl bromide          |                                   | C <sub>2</sub> Br <sub>2</sub> O <sub>2</sub> | 15219-34-8 | 10.49     | 5    |    |      |
| 183 | Oxydiethanol, 2,2-      | diethylene glycol                 | C <sub>4</sub> H <sub>10</sub> O <sub>3</sub> | 111-46-6   | ~10.<br>3 | 2    | 20 | 400  |
| 184 | Oxygen                  |                                   | 02  | 7782-44-7  | 12.07     | NR   |    |      |
| 185 | Ozone                   |                                   | O <sub>3</sub>                                | 10028-15-6 | 12.52     | NR   |    |      |
| 186 | Paraffins, normal       |                                   |   | 64771-72-8 | ~9.5      | 1    | 5  | 100  |
| 187 | Paraldehyde             |                                   | C <sub>6</sub> H <sub>12</sub> O <sub>3</sub> | 123-63-7   | ~9.7      | 2.2  |    |      |
| 188 | Pentane                 |                                   | C <sub>5</sub> H <sub>12</sub>                | 109-66-0   | 10.35     | 7    | 40 | 800  |
| 189 | Pentanoic acid          |                                   | C <sub>5</sub> H <sub>10</sub> O <sub>2</sub> | 109-52-4   | 10.53     | 8    |    |      |
| 190 | Pentanol, 2-            |                                   | C <sub>5</sub> H <sub>12</sub> O              | 6032-29-7  | 9.78      | 2    |    |      |
| 191 | Pentanol, 3-            |                                   | C <sub>5</sub> H <sub>12</sub> O              | 584-02-1   | 9.76      | 1.7  |    |      |
| 192 | Pentene, 1-             |                                   | C <sub>5</sub> H <sub>10</sub>                | 109-67-1   | 9.49      | 0.92 |    |      |
| 193 | Pentyne, 1-             |                                   | C <sub>5</sub> H <sub>8</sub>                 | 627-19-0   | 10.1      | 3    |    |      |
| 194 | Peracetic acid          |                                   | C <sub>2</sub> H <sub>4</sub> O <sub>3</sub>  | 79-21-0    | ~10.<br>5 | 2    | 10 | 200  |
| 195 | Petroleum ether         | ligroin, VM&P<br>naphtha, benzine |   | 8032-32-4  | ~10       | 0.9  |    |      |
| 196 | Phenol                  | hydroxybenzene                    | C <sub>6</sub> H <sub>6</sub> O               | 108-95-2   | 8.51      | 1.2  | 6  | 120  |
| 197 | Phosgene                |                                   | COCI <sub>2</sub>                             | 75-44-5    | 11.55     | NR   |    |      |
| 198 | Phosphine               |                                   | PH <sub>3</sub>                               | 7803-51-2  | 9.96      | 2    | 10 | 200  |
| 199 | Propane-1,2-diol        | propylene glycol                  | C <sub>3</sub> H <sub>8</sub> O <sub>2</sub>  | 57-55-6    | 10        | 3    | 50 | 1000 |
| 200 | Acetaldehyde            |                                   | C <sub>3</sub> H <sub>9</sub> NO              | 156-87-6   | ~9.5      | 1.5  |    |      |
| 201 | Propene                 | propylene                         | C <sub>3</sub> H <sub>6</sub>                 | 115-07-1   | 9.73      | 1.4  | 7  | 140  |

| 202   | Propiolic acid                       | 2-propynoic acid  | C <sub>3</sub> H <sub>2</sub> O <sub>2</sub>   | 471-25-0   | 10.45     | 8                                  |                           |                            |
|-------|--------------------------------------|---|--|------------|-----------|------------------------------------|---------------------------|----------------------------|
| 203   | Propionaldehyde                      | propanal,<br>propional                                  | C <sub>3</sub> H <sub>6</sub> O                | 123-38-6   | 9.95      | 1.7                                | 8                         | 169                        |
| Index | Chemical name                        | Alternative name  | Formula  | CAS no.    | IE, eV    | Response<br>Factor (RF)<br>10.6 eV | E2618-<br>PID-40<br>(ppb) | E2618-<br>PID-200<br>(ppb) |
| 204   | Propionic acid                       |   | C <sub>3</sub> H <sub>6</sub> O <sub>2</sub>   | 79-09-4    | 10.44     | 8                                  | 40                        | 800                        |
| 205   | Propyl acetate, n-                   |   | C <sub>5</sub> H <sub>10</sub> O <sub>2</sub>  | 109-60-4   | 10.04     | 3                                  | 13                        | 250                        |
| 206   | Propyl iodide                        | lodopropane   | C <sub>3</sub> H <sub>7</sub> I                | 107-08-4   | 9.26      | 1                                  |                           |                            |
| 207   | Propylamine, n-                      |   | C <sub>3</sub> H <sub>9</sub> N                | 107-10-8   | 8.5       | 1.1                                |                           |                            |
| 208   | Propylbenzene (all isomers)          |   | C <sub>9</sub> H <sub>12</sub>                 | 74296-31-4 | 8.7       | 0.5                                |                           |                            |
| 209   | Propylene carbonate                  |   | C <sub>4</sub> H <sub>6</sub> O <sub>3</sub>   | 108-32-7   | ~10.<br>5 | 15                                 |                           |                            |
| 210   | Propylene dinitrate                  |   | C <sub>3</sub> H <sub>6</sub> N <sub>2</sub> O | 6423-43-4  | ~11       | NR                                 |                           |                            |
| 211   | Propylene glycol ethyl ether acetate | PGEEA   | C <sub>7</sub> H <sub>14</sub> O <sub>3</sub>  | 98516-30-4 | ~9.6      | 1.2                                |                           |                            |
| 212   | Propylene oxide                      |   | C <sub>3</sub> H <sub>6</sub> O                | 75-56-9    | 10.22     | 6                                  | 35                        | 700                        |
| 213   | Propyleneimine                       | 2-methylaziridine                                       | C <sub>3</sub> H <sub>7</sub> N                | 75-55-8    | 9         | 1.4                                | 7                         | 130                        |
| 214   | Propyl Nitrate, n-                   |   | C <sub>3</sub> H <sub>7</sub> NO <sub>3</sub>  | 627-13-4   | 11.07     | NR                                 |                           |                            |
| 215   | Propyne                              | methylacetylene   | C <sub>3</sub> H <sub>4</sub>                  | 74-99-7    | 10.36     | 4                                  |                           |                            |
| 216   | Pyrazine                             |   | C <sub>4</sub> H <sub>4</sub> N <sub>2</sub>   | 290-37-9   | 9.29      | 3                                  |                           |                            |
| 217   | Pyridine                             |   | C <sub>5</sub> H <sub>5</sub> N                | 110-86-1   | 9.25      | 0.7                                | 4                         | 75                         |
| 218   | Pyrrole                              |   | C₄H₅N  | 109-97-7   | 8.02      | 0.6                                |                           |                            |
| 219   | Pyrrolidine                          |   | C <sub>4</sub> H <sub>9</sub> N                | 123-75-1   | 8.77      | 4                                  |                           |                            |
| 220   | Pyruvaldehyde                        |   | C <sub>3</sub> H <sub>4</sub> O <sub>2</sub>   | 78-98-8    | 9.6       | 0.7                                |                           |                            |
| 221   | Rose oxide, cis-                     | Tetrahydro-4-met<br>hyl-2-(2-methylpr<br>opyl)-2H-pyran | C <sub>10</sub> H <sub>18</sub> O              | 16409-43-1 | ~9        | 0.8                                |                           |                            |
| 222   | Sec-amyl acetate                     |   | C <sub>7</sub> H <sub>14</sub> O <sub>2</sub>  | 626-38-0   | ~9.9      | 5                                  |                           |                            |
| 223   | Sevoflurane                          | 1,1,1,3,3,3-hexafl<br>uoro-2-(fluorome<br>thoxy)propane | C₃H₃F <sub>7</sub> O                           | 28523-86-6 | 11        | NR                                 |                           |                            |
| 224   | Styrene                              | vinylbenzene  | C <sub>8</sub> H <sub>8</sub>                  | 100-42-5   | 8.4       | 0.45                               | 2                         | 50                         |
| 225   | Tetrafluoroethylene                  |   | C <sub>2</sub> F <sub>4</sub>                  | 116-14-3   | 10.12     | 15                                 | 5                         | 100                        |
| 226   | Tetrafluoromethane                   | carbon<br>tetrafluoride                                 | CF₄  | 75-73-0    | 15.3      | NR                                 |                           |                            |
| 227   | Tetrahydrofuran                      | THF   | C <sub>4</sub> H <sub>8</sub> O                | 109-99-9   | 9.41      | 2.3                                | 8                         | 150                        |
| 228   | Toluene                              |   | C <sub>7</sub> H <sub>8</sub>                  | 108-88-3   | 8.82      | 0.56                               | 3                         | 50                         |
| 229   | Triethylamine                        | TEA   | C <sub>6</sub> H <sub>15</sub> N               | 121-44-8   | 7.5       | 1.3                                | 5                         | 90                         |

| 230 | Trifluoroacetic acid | TFAA                   | C <sub>2</sub> HO <sub>2</sub> F <sub>3</sub> | 76-05-1   | 11.46 | NR   |   |     |
|-----|----------------------|------------------------|---|-----------|-------|------|---|-----|
| 231 | Trifluoroethane      | trifluoroethylene      | C <sub>2</sub> HF <sub>3</sub>                | 359-11-5  | 10.14 | 5    |   |     |
| 232 | Turpentine           |                        | C <sub>10</sub> H <sub>16</sub>               | 9005-90-7 | ~8.5  | 0.6  |   |     |
| 233 | Turpentine oil       | pinenes                | C <sub>10</sub> H <sub>16</sub>               | 8006-64-2 | ~8    | 0.6  | 3 | 60  |
| 234 | Undecane             |                        | C <sub>11</sub> H <sub>24</sub>               | 1120-21-4 | 9.56  | 1.1  | 5 | 100 |
| 235 | Vanillin             |                        | C <sub>8</sub> H <sub>8</sub> O <sub>3</sub>  | 121-33-5  | ~9    | 1    |   |     |
| 236 | Water                | dihydrogen<br>monoxide | H <sub>2</sub> O                              | 7732-18-5 | 12.61 | NR   |   |     |
| 237 | Xenon                |                        | Xe  | 7440-63-3 | 12.13 | NR   |   |     |
| 238 | Xylene mixed isomers | dimethyl<br>benzenes   | C <sub>8</sub> H <sub>10</sub>                | 1330-20-7 | 8.56  | 0.54 | 2 | 40  |
| 239 | Xylene, m-           |                        | C <sub>8</sub> H <sub>10</sub>                | 108-38-3  | 8.56  | 0.5  | 2 | 50  |
| 240 | Xylene, o-           |                        | C <sub>8</sub> H <sub>10</sub>                | 95-47-6   | 8.56  | 0.5  | 3 | 60  |
| 241 | Xylene, p-           |                        | C <sub>8</sub> H <sub>10</sub>                | 106-42-3  | 8.44  | 0.55 | 3 | 50  |

### **Calibration and maintenance**

Do not perform any maintenance operation with the power on.

Clean the device with a soft damp cloth. Do not use any abrasive cleaning agents. Do not immerse the device in water or any cleaning media.

**NOTE!** Cleaning instruments and spare parts are not included in the delivery set, Please contact your Seller for more information.

#### How often the PID needs maintenance?

The frequency of calibration and maintenance depends on operating conditions (VOC concentration, humidity, dust, and oil in the air, etc).

In clean indoor air with low VOC concentration, a monthly or less frequent calibration is recommended. When higher VOC concentrations are measured under harsh conditions, perform calibration and maintenance more frequently.

#### PID needs maintenance if:

- The baseline is climbing after you zero the PID replace the electrode stack.
- The PID becomes sensitive to humidity replace electrode stack.
- The PID cell shows signs of contamination after the lamp window has been cleaned – replace the electrode stack.
- The PID cell is known to have been subjected to severe contamination—replace the electrode stack
- The baseline is unstable or shifts when you move the PID replace the electrode stack
- The sensitivity has dropped too much (note the change required when checking
- calibration) clean the lamp
- The signal to noise ratio at low VOC concentrations becomes inadequate replace the lamp

Cleaning of the PID lamp is recommended as the first action when a PID needs maintenance. It is recommended to recalibrate the device every time after stack replacement or lamp cleaning.

#### Removing the electrode stack and lamp

**NOTE!** Always use the Electrode Stack Removal Tool to remove the electrode stack; any other tools may damage your PID and invalidate your warranty.

To remove the electrode stack, proceed as follows:

- 1. Gently remove the sensor from the equipment.
- 2. Place the PID, pellet side down, onto a clean surface.
- 3. Locate the electrode stack removal tool into the two slots on the sides of the PID and squeeze together until the electrode stack and lamp are released.
- 4. Carefully lift the PID body away from the pellet and lamp.
- Occasionally the lamp may be temporarily lodged in the cell and will need to be freed carefully with tweezers.
- 6. Occasionally the small spring behind the lamp will come out when the lamp is removed from the sensor. Simply place it into the sensor house.



Electrode stack removing tool



Electrode stack

#### Cleaning the PID Lamp

Inspection of the lamp may reveal a layer of contamination on the detection window that presents itself as a 'blue hue.' To check for contamination, hold the lamp in front of a light source and look across the window surface. To avoid contaminating the sensor and affecting accuracy, do not touch the lamp window with bare fingers. You may touch the body of the lamp with clean fingers.

**NOTE!** Always use the recommended cleaning kit to clean the lamp.

#### PID lamp cleaning kit

The vial of the cleaning compound contains alumina (CAS Number 1344-28-1) as a very fine powder. Please contact your Seller for a full material safety data sheet MSDS.

Key safety issues are identified below:

Hazard identification: May cause irritation of the respiratory tract and eyes.

Storage: Keep container closed to prevent water absorption and contamination

#### Handling:

- Do not breathe in the powder. Avoid contact with skin, eyes, and clothing;
- Wear suitable protective clothing;
- Follow industrial hygiene practices: Wash face and hands thoroughly with soap and water after use and before eating, drinking, smoking, or applying cosmetics.
- The powder carries a TVL (TWA) limit of 10 mg/m<sup>3</sup>

#### PID lamp cleaning instructions:

- 1. Open the container of the alumina polishing compound;
- 2. With a clean cotton bud, collect a small amount of the powder;
- Use this cotton bud to polish the PID lamp window. Use a circular action, applying light pressure to clean the lamp window. Do not touch the lamp window with fingers;
- Continue polishing until an audible "squeaking" is made by the cotton bud moving over the window surface (usually within 15 seconds);
- 5. Remove the residual powder from the lamp window with a clean cotton bud. Care must be taken not to touch the tips of cotton buds that are to be used to clean the lamps as this may contaminate them with finger oil.
- 6. Ensure the lamp is completely dry and remove any visible signs of contamination before refitting.

#### Re-fitting the PID electrode stack and lamp

**NOTE!** Never refit a damaged lamp.

 Place the lamp inside the O-ring seal in the pellet as illustrated below. Twisting the lamp slightly during insertion will help to ensure the lamp window is snug against the electrode stack's front electrode. The lamp should be freely supported by the O-ring.

- Continuing to hold the electrode stack between forefinger and thumb, carefully insert the lamp into the recess in the sensor ensuring that the lamp remains in position. Press the electrode stack firmly, to ensure that the electrode stack wing clips are engaged, and the top faces of the electrode stack and sensor house are flush.
- 3. Refit the sensor into the sensing equipment.
- 4. Recalibrate the gas detector.



#### Discarding the PID electrode stack

The electrode stack does not have any toxic components, however, if it has been contaminated by toxic materials, show due care when disposing of.

# **Delivery set**

• Gas transmitter E2618 (wall mount, duct mount, or remote probe version)

- Mounting accessories:
  - 4 screws with plastic dowels for wall mount version
  - Square Rubber flange for Duct mount option
  - o Round rubber flange for Remote probe option
  - o Fixing clamp for Remote probe option







### Order code for E2618-PID options

| E2618 options  | Order code        |
|--|-------------------|
| Remote probe, 3 m cable  | E2618-PID-RP33-3  |
| Remote probe, 10 m cable   | E2618-PID-RP33-10 |
| Duct mount option, stem Ø35×L230 mm  | E2618-PID-DM      |
| Integrated 24 VAC power supply module                                      | E2618-PID-24VAC   |
| Remote LCD single value display, wall mount box 115 × 65 × 40mm, 3 m cable | E2618-PID-RLCD3   |

# Configuring

A standard configuration kit includes a USB-RS485 converter, fixed flow regulator, gas tubing with applicators, and a software pack. Please contact your Seller for more information.

Gas transmitters E2618 share all functionalities of the PluraSens® multifunctional transmitter platform. The features and options include:

- Digital output change rate-limiting filter
- Digital integrating (averaging) filter
- Free assignment of each analog output to the chosen parameter
- Flexible setting of analog output scales for each output
- Output zero and slope adjustment for calibration

The output scales can be changed by Modbus commands by using the configuration software and the standard configuration kit (see Modbus RTU Communication).

# Return to default settings

To reset the device's Slave ID, baud rate, and stop bit numbers to factory settings, proceed as follows:

- 1. De-energize the device
- 2. Connect the X6 jumper
- 3. Turn on the device
- 4. De-energize the device
- 5. Disconnect the X6 jumper
- 6. Turn on the device

### **Modbus RTU Communication**

#### RS485 communication interface

| Databits: 8 Parity: none / odd / even Stop bits: 1 or 2 Protocol: Modbus RTU | Supported Modbus functions:<br>03 – Read multiple registers<br>06 – Write a single register |
|--|---|
|--|---|

### **Communication parameters**

| Parameter            | Permitted values  | Default |
|----------------------|---|---------|
| Supported baud rates | 1200, 2400, 4800, 9600, 19200, 38400, 57600   | 9600    |
| Data bits            | 8   | 8       |
| Parity               | none / odd / even   | none    |
| Stop bits            | 1, 2  | 1       |
| Protocol             | Modbus RTU  |         |
| Modbus functions     | 03 - Read multiple registers<br>06 - Write a single register  |         |
| Error codes          | 01 – Illegal function 02 – Illegal data address 03 – Illegal data value 04 – Slave device failure (details of last error 04 can be read from register 0x000 | 08)     |

# **Modbus holding registers**

Register addresses are shown 0-based, Address in hexadecimal, Reg in decimal format. Modbus holding register numbers MHR are shown in the decimal 1-based format and may be addressed either from 00001 or 40001 base.

| Address | Reg / MHR | RW | Description              | Supported values (dec) | Default |
|---------|-----------|----|--------------------------|------------------------|---------|
| 0x0001  | 1 / 40002 | R  | Hardware version         |                        | -       |
| 0x0002  | 2 / 40003 | R  | Software version         |                        | -       |
| 0x0003  | 3 / 40004 | R  | Product serial number    | 165535                 | -       |
| 0x0004  | 4 / 40005 | RW | Slave ID (net address) * | 1247 **                | 1       |

| 0x0005  | 5 / 40006      | RW | Baud rate *   | 1200, 2400, 4800, 9600, 19200,<br>38400, 57600   | 9600    |
|---------|----------------|----|---|--|---------|
| 0x0006  | 6 / 40007      | RW | Response delay, ms  | 1255   | 10      |
| 0x0007  | 7 / 40008      | RW | Stop bits, parity bit *   | 1 - No parity bit, 1 stop bit<br>(default after factory reset)<br>2 - No parity bit, 2 stop bits<br>3 - Odd parity, 1 stop bit<br>4 - Even parity, 1 stop bit<br>NOTE: 3 and 4 are available<br>starting from the Software<br>version 0x218 (dec. 536) | 1       |
| 0x0008  | 8 /40009       | R  | Last error code   | 1255   | -       |
| 0x0011  | 17 / 40018     | RW | Technological:<br>age of last data in seconds<br>(read) / restart (write) | 065535 s (read),<br>writing 42330 restarts the<br>instrument   | -       |
|         |                |    |   |  |         |
| 0x00A5  | 165 /<br>40166 | RW | Zero adjustments for gas data, ADC  | -32000+32000 ADC units   | 0       |
| 0x00A6  | 166 /<br>40167 | RW | Slope adjustment for gas data   | 165535   | 512     |
| 0x00A7  | 167 /<br>40168 | RW | Change rate limit for gas<br>data, ppm (% for O2) / s                     | 132000, 0 - no limit   | 0       |
| 0x00A8  | 168 /<br>40169 | RW | Integrating filter time constant, s                                       | 132000 (seconds),<br>0 - no filter   | 0       |
| Address | Reg / MHR      | RW | Description   | Supported values (dec)   | Default |
| 0x00C9  | 201 /<br>40202 | RW | Parameter tied to analog output 1   | 0 - None<br>2 - Gas concentration<br>9 - Forced Modbus control,<br>value set in MHR / 40204  | 2       |
| 0x00CA  | 202 /<br>40203 | RW | Parameter tied to analog output 2   | 0 - None<br>2 - Gas concentration<br>9 - Forced Modbus control,<br>value set in MHR / 40205  | 2       |
| 0x00CB  | 203 /<br>40204 | RW | Forced value for analog output 1***                                       | 01000<br>(0.0%100.0% of output scale)  | 0       |
| 0x00CC  | 204 /<br>40205 | RW | Forced value for analog output 2***                                       | 01000<br>(0.0%100.0% of output scale)  | 0       |
|         |                |    |   |  | _       |

| 0x00FF  | 255 /<br>40256 | RW | Sensor, analog outputs,<br>LED and buzzer status | bit[0]=0/1 - sensor present/absent, RO bit[1]=0/1 - analog outputs deactivated/activated bit[2]=0/1 - in case the sensor is absent, turn signaling off/on analog output1 bit[3]=0/1 - in case the sensor is absent, turn on signaling with low current/high current on analog output1; if bit[2]=0 this bit will be ignored bit[4]=0/1 - in case of sensor absent, turn signaling off/on analog output2 bit[5]=0/1 - in case of sensor absent, turn on signaling with low current/high current on analog output2 bit[5]=0/1 - in case of sensor absent, turn on signaling with low current/high current on analog output2; if bit[4]==0 this bit will be ignored bit[6]=0/1 - current/voltage output detected on output1, RO bit[7]=0/1 - current/voltage output detected on output2, RO bit[8]=0/1 - LED deactivated/activated (always 0 for E2618) bit[10]=0/1 - LED is on/off in normal condition |                 |
|---------|----------------|----|--|--|-----------------|
| Address | Reg / MHR      | RW | Description                                      | Supported values (dec)   | Default         |
| 0x0101  | 257 /<br>40258 | R  | Raw gas sensor data                              | 04095, ADC units   |                 |
| 0x0103  | 259 /<br>40260 | R  | Gas concentration, gas units                     | 065535, gas units  |                 |
| 0x0105  | 261 /<br>40262 | RW | 0% value<br>for analog output 1                  | -32000+32000<br>(gas unit/°C)  | User<br>defined |
| 0x0106  | 262 /<br>40263 | RW | 100% value<br>for analog output 1                | -32000+32000<br>(gas unit/°C)  | User<br>defined |
| 0x0107  | 263 /<br>40264 | RW | 0% value<br>for analog output 2                  | -32000+32000<br>(gas unit/°C)  | User<br>defined |
| 0x0108  | 264 /<br>40265 | RW | 100% value<br>for analog output 2                | -32000+32000<br>(gas unit/°C)  | User<br>defined |

- \* The new value is applied after restart.
- \*\* Broadcast slave ID 0 can be used to assign a new ID to the instrument with an unknown ID. When addressing by ID 0 the device shall be the only Modbus instrument in the network. The device will not respond to the Master command when addressed by ID 0.

\*\*\* - This value is dynamic and not kept in EEPROM after a restart.

# Warranty

This product is warranted to be free from defects in material and workmanship for a period of one year from the date of the original sale. During this warranty period, the Manufacturer will, at its option, either repair or replace a product that proves to be defective. This warranty is void if the product has been operated in conditions outside ranges specified by the Manufacturer or damaged by customer error or negligence or if there has been an unauthorized modification.

### Manufacturer contacts

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