

MiniPID HI or MiniPID LO photoionisation cells are calibrated using isobutylene, but PID is a broadband VOC detection technique, 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. Remember, these are approximate values, so for best accuracy you should calibrate with the relevant VOC.

The table includes six columns:

- 1 **Gas/ VOC** The most common name for the VOC. If you can not find the name of your VOC of concern, then email us at info@ionscience.com and we will help.
- 2 **CAS No.** Sometimes it is easier to identify a VOC from the internationally recognised CAS No.: ask your supplier.
- 3 **Formula** To further assist in identifying the VOC, this also is helpful in identifying its molecular weight, from which ppm measurements can be converted to say, mg/m³ measurements.
- 4 **Response Factor (RF)** also known as correction factor. Multiply the output response from the cell by the RF to provide a normalised scale of VOC concentration.
- 5 **Relative sensitivity (%)** This 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 the same way as cross-sensitivity for toxic gas sensors.
- 6 **Minimum Detection Level (MDL) or Minimum Detectable Quantity (MDQ).** Typical lowest concentration that can be detected. The Mini PID HI has greater sensitivity than the Mini PID LO, so MDL's for the Mini PID HI are much lower.

The RF is measured in dry air; high humidity will reduce this factor by 10% to 20%, so the RF should be increased in high humidities.

VOC response

The PID can not measure all VOCs or gases. Two types of VOCs are not measured:

ZR: No response. The 10.6 eV lamp does not ionise the VOC and the VOC can not be measured.

NV: The vapour pressure of the VOC at 20°C is less than a few ppm, so this Semi-Volatile Organic Compound can not be measured.

Frequently 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 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, balance nitrogen. Some gases absorb UV light (oxygen, methane), so in gases where there are significant concentrations of oxygen or methane, the apparent concentration will be less than is actually present. Methane absorbs UV strongly, so for accurate measurements in methane, calibrate with your target VOC in the expected methane concentration. 50% LEL methane reduces reading by up to 50%. Gases such as nitrogen and helium do not absorb UV and do not affect the relative response.

Accuracy of the Table

This table is for indication only. The values are only accurate within the linear range of the sensor (to 300 ppm), although this linear range may vary somewhat for specific gases. Table accuracy is to 1 or 2 digits only, so when calculating concentration for a specific VOC, specify to 1 or 2 digits only. For best accuracy, calibrate using the specific VOC, as advised in the 'MiniPID user guide'.

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| Gas/ VOC | CAS No. | Formula | Relative response | Relative sensitivity (%) | Typical MDL Mini PID HI (ppb) | Typical MDL Mini PID LO (ppb) |
|-----------------------------|-----------|---------|-------------------|--------------------------|-------------------------------|-------------------------------|
| Acetaldehyde | 75-07-0 | C2H4O | 4.9 | 21 | 25 | 480 |
| Acetic Acid | 64-17-7 | C2H4O2 | 36 | 3 | 180 | 3615 |
| Acetic Anhydride | 108-24-7 | C4H6O3 | 4.0 | 25 | 20 | 400 |
| Acetone | 67-64-1 | C3H6O | 0.7 | 140 | 5 | 70 |
| Acetonitrile | 75-05-8 | CH3CN | ZR | | | |
| Acetylene | 74-86-2 | C2H2 | ZR | | | |
| Acrolein | 107-02-8 | C3H4O | 4.0 | 25 | 20 | 400 |
| Acrylic Acid | 79-10-7 | C3H4O2 | 2.7 | 36 | 15 | 275 |
| Acrylonitrile | 107-13-1 | C3H3N | ZR | | | |
| Allyl alcohol | 107-18-6 | C3H6O | 2.1 | 48 | 10 | 200 |
| Allyl chloride | 107-05-1 | C3H5Cl | 4.5 | 22 | 20 | 450 |
| Ammonia | 7664-41-7 | H3N | 8.5 | 12 | 40 | 850 |
| Amyl acetate, n- | 628-63-7 | C7H14O2 | 1.8 | 56 | 10 | 180 |
| Amyl alcohol | 71-41-0 | C5H12O | 3.2 | 31 | 15 | 320 |
| Aniline | 62-53-3 | C6H7N | 0.5 | 200 | 3 | 50 |
| Anisole | 100-66-3 | C7H8O | 0.5 | 211 | 2 | 50 |
| Arsine | 7784-42-1 | AsH3 | 2.5 | 40 | 15 | 250 |
| Asphalt, petroleum fumes | 8052-42-4 | | 1.0 | 100 | 5 | 100 |
| Benzaldehyde | 100-52-7 | C7H6O | 0.9 | 117 | 5 | 85 |
| Benzene | 71-43-2 | C6H6 | 0.5 | 200 | 3 | 50 |
| Benzenethiol | 108-98-5 | C6H5SH | 0.7 | 143 | 4 | 70 |
| Benzonitrile | 100-47-0 | C7H5N | 0.7 | 141 | 4 | 70 |
| Benzyl alcohol | 100-51-6 | C7H8O | 1.3 | 80 | 6 | 125 |
| Benzyl chloride | 100-44-7 | C7H7Cl | 0.6 | 182 | 3 | 55 |
| Benzyl formate | 104-57-4 | C8H8O2 | 0.8 | 130 | 5 | 77 |
| Biphenyl | 92-52-4 | C12H10 | 0.4 | 250 | 2 | 40 |
| Bis(2,3-epoxypropyl) ether | 2238-07-5 | C6H10O3 | 3.0 | 33 | 15 | 300 |
| Boron trifluoride | 7637-07-2 | BF3 | ZR | | | |
| Bromine | 7726-95-6 | Br2 | 20 | 5 | 100 | 2000 |
| Bromine pentafluoride | 7789-30-2 | BrF5 | ZR | | | |
| Bromobenzene | 108-86-1 | C6H5Br | 0.7 | 143 | 4 | 70 |
| Bromochloromethane | 74-97-5 | CH2ClBr | ZR | | | |
| Bromoethane | 74-96-4 | C2H5Br | 5.0 | 20 | 25 | 500 |
| Bromoethyl methyl ether, 2- | 6482-24-2 | C3H7OBr | 2.5 | 40 | 15 | 250 |
| Bromoform | 75-25-2 | CHBr3 | 2.8 | 36 | 15 | 280 |
| Bromopropane, 1- | 106-94-5 | C3H7Br | 1.3 | 77 | 7 | 130 |
| Bromotrifluoromethane | 75-63-8 | CF3Br | ZR | | | |
| Butadiene | 106-99-0 | C4H6 | 0.8 | 120 | 4 | 80 |
| Butadiene diepoxyde, 1,3- | 1464-53-5 | C4H6O2 | 4.0 | 25 | 20 | 400 |
| Butane, n- | 106-97-8 | C4H10 | 46 | 2 | 230 | 4600 |
| Butanol, 1- | 71-36-3 | C4H10O | 4.0 | 25 | 20 | 400 |
| Buten-3-ol, 1- | 598-32-3 | C4H8O | 1.2 | 87 | 6 | 115 |
| Butene, 1- | 106-98-9 | C4H8 | 1.3 | 77 | 7 | 130 |
| Butoxyethanol, 2- | 111-76-2 | C6H14O2 | 1.1 | 91 | 6 | 110 |
| Butyl acetate, n- | 123-86-4 | C6H12O2 | 2.4 | 41 | 10 | 240 |
| Butyl acrylate, n- | 141-32-2 | C7H12O2 | 1.5 | 67 | 8 | 150 |
| Butyl lactate | 138-22-7 | C7H14O3 | 2.5 | 40 | 15 | 250 |
| Butyl mercaptan | 109-79-5 | C4H10S | 0.5 | 185 | 3 | 50 |
| Butylamine, 2- | 513-49-5 | C4H11N | 0.9 | 111 | 5 | 90 |
| Butylamine, n- | 109-73-9 | C4H11N | 1.0 | 100 | 5 | 100 |
| Camphene | 565-00-4 | C10H16 | 0.5 | 222 | 2 | 45 |
| Carbon dioxide | 124-38-9 | CO2 | ZR | | | |

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| Gas/ VOC | CAS No. | Formula | Relative response | Relative sensitivity (%) | Typical MDL Mini PID HI (ppb) | Typical MDL Mini PID LO (ppb) |
|----------------------------------|------------|----------|-------------------|--------------------------|-------------------------------|-------------------------------|
| Carbon disulfide | 75-15-0 | CS2 | 1.4 | 71 | 7 | 140 |
| Carbon monoxide | 630-08-0 | CO | ZR | | | |
| Carbon tetrabromide | 558-13-4 | CBr4 | 3.0 | 33 | 15 | 300 |
| Carbon tetrachloride | 56-23-5 | CCl4 | ZR | | | |
| Carbonyl sulphide | 463-58-1 | COS | ZR | | | |
| Carvone, R- | 6485-40-1 | C10H14O | 1.0 | 100 | 5 | 100 |
| Chlorine | 7782-50-5 | Cl2 | ZR | | | |
| Chlorine dioxide | 10049-04-4 | ClO2 | 1.0 | 100 | 5 | 100 |
| Chlorine trifluoride | 7790-91-2 | ClF3 | ZR | | | |
| Chloro-1,1,1,2-tetrafluoroethane | 2837-89-0 | C2HClF4 | ZR | | | |
| Chloro-1,1,1-trifluoroethane, 2- | 75-88-7 | C2H2ClF3 | ZR | | | |
| Chloro-1,1,2,2-tetrafluoroethane | 354-25-6 | C2HClF4 | ZR | | | |
| Chloro-1,1,2-trifluoroethane, 1- | 421-04-5 | C2H2ClF3 | ZR | | | |
| Chloro-1,1-difluoroethane, 1- | 75-68-3 | C2H3ClF2 | ZR | | | |
| Chloro-1,1-difluoroethane, 1- | 75-68-3 | C2H3ClF2 | ZR | | | |
| Chloro-1,1-difluoroethane, 2- | 338-65-8 | C2H3ClF2 | ZR | | | |
| Chloro-1,2,2-trifluoroethane | 431-07-2 | C2H2ClF3 | ZR | | | |
| Chloro-1,3-butadiene, 2- | 126-99-8 | C4H5Cl | 3.2 | 30 | 16 | 320 |
| Chloro-1-fluoroethane, 1- | 1615-75-4 | C2H4ClF | ZR | | | |
| Chloro-2-fluoroethane, 1- | 762-50-5 | C2H4ClF | ZR | | | |
| Chloroacetaldehyde | 107-20-0 | C2H3OCl | ZR | | | |
| Chlorobenzene | 108-90-7 | C6H5Cl | 0.5 | 220 | 2 | 50 |
| Chlorodifluoromethane | 75-45-6 | CHClF2 | ZR | | | |
| Chloroethane | 75-00-3 | C2H5Cl | ZR | | | |
| Chloroethanol 2- | 107-07-3 | C2H5ClO | 10.0 | 10 | 50 | 1000 |
| Chloroethyl methyl ether, 2- | 627-42-9 | C3H7ClO | 2.6 | 40 | 13 | 250 |
| Chlorofluoromethane | 593-70-4 | CH2ClF | ZR | | | |
| Chloroform | 67-66-3 | CHCl3 | ZR | | | |
| Chloromethane | 74-87-3 | CH3Cl | ZR | | | |
| Chloropentafluoroethane | 76-15-3 | C2ClF5 | ZR | | | |
| Chlorotoluene, o- | 95-49-8 | C7H7Cl | 0.5 | 220 | 2 | 50 |
| Chlorotoluene, p- | 108-41-8 | C7H7Cl | 0.5 | 200 | 3 | 50 |
| Chlorotrifluoroethylene | 79-38-9 | C2ClF3 | 1.0 | 100 | 5 | 100 |
| Chlorotrifluoromethane | 75-72-9 | CClF3 | ZR | | | |
| Citral | 5392-40-5 | C10H16O | 1.0 | 100 | 5 | 100 |
| Citronellol | 26489-01-0 | C10H20O | 1.0 | 100 | 5 | 100 |
| Cresol, m- | 108-39-4 | C7H8O | 1.1 | 95 | 5 | 105 |
| Cresol, o- | 95-48-7 | C7H8O | 1.1 | 95 | 5 | 105 |
| Cresol, p- | 106-44-5 | C7H8O | 1.1 | 95 | 5 | 105 |
| Crotonaldehyde | 4170-30-3 | C4H6O | 1.0 | 100 | 5 | 100 |
| Cumene | 98-82-8 | C9H12 | 0.6 | 170 | 3 | 60 |
| Cyanamide | 420-04-2 | CH2N2 | ZR | | | |
| Cyanogen bromide | 506-68-3 | CNBr | ZR | | | |
| Cyanogen chloride | 506-77-4 | CNCI | ZR | | | |
| Cyclohexane | 110-82-7 | C6H12 | 1.3 | 77 | 7 | 130 |
| Cyclohexanol | 108-93-0 | C6H12O | 2.9 | 34 | 15 | 300 |
| Cyclohexanone | 108-94-1 | C6H10O | 1.1 | 91 | 6 | 110 |
| Cyclohexene | 110-83-8 | C6H10 | 0.8 | 133 | 5 | 75 |
| Cyclohexylamine | 108-91-8 | C6H13N | 1.0 | 102 | 5 | 100 |
| Cyclopentane | 287-92-3 | C5H10 | 4.0 | 25 | 20 | 400 |
| Decane, n- | 124-18-5 | C10H22 | 1.0 | 96 | 5 | 100 |
| Diacetone alcohol | 123-42-2 | C6H12O2 | 0.8 | 125 | 5 | 80 |
| Dibenzoyl peroxide | 94-36-0 | C14H10O4 | 0.8 | 125 | 5 | 80 |

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| Gas/ VOC | CAS No. | Formula | Relative response | Relative sensitivity (%) | Typical MDL Mini PID HI (ppb) | Typical MDL Mini PID LO (ppb) |
|--------------------------------------|------------|-----------|-------------------|--------------------------|-------------------------------|-------------------------------|
| Diborane | 19287-45-7 | B2H6 | ZR | | | |
| Dibromochloromethane | 124-48-1 | CHBr2Cl | 10.0 | 10 | 50 | 1000 |
| Dibromodifluoromethane | 75-61-6 | CF2Br2 | ZR | | | |
| Dibromoethane, 1,2- | 106-93-4 | C2H4Br2 | 2.0 | 50 | 10 | 200 |
| Dibromotetrafluoroethane, 1,2- | 124-73-2 | C2F4Br2 | ZR | | | |
| Diethyl hydrogen phosphate | 107-66-4 | HC8H18PO4 | 4.0 | 25 | 20 | 400 |
| Dichloro-1,1,1-trifluoroethane, 2,2- | 306-83-2 | C2HCl2F3 | ZR | | | |
| Dichloro-1,1-difluoroethane, 1,2- | 1649-08-7 | C2H2Cl2F2 | ZR | | | |
| Dichloro-1,2,2-trifluoroethane, 1,2- | 354-23-4 | C2HCl2F3 | ZR | | | |
| Dichloro-1,2-difluoroethane, 1,2- | 631-06-1 | C2H2Cl2F2 | ZR | | | |
| Dichloro-1-fluoroethane, 1,1- | 1717-00-6 | C2H3Cl2F | ZR | | | |
| Dichloro-1-fluoroethane, 1,1- | 1717-00-6 | C2H3Cl2F | ZR | | | |
| Dichloro-1-fluoroethane, 1,2- | 430-57-9 | C2H3Cl2F | ZR | | | |
| Dichloro-1-propene, 2,3- | 78-88-6 | C3H4Cl2 | 1.4 | 70 | 7 | 140 |
| Dichloro-2,2-difluoroethane, 1,1- | 79-35-6 | C2H2Cl2F2 | ZR | | | |
| Dichloroacetylene | 7572-29-4 | C2Cl2 | 5.0 | 20 | 25 | 500 |
| Dichlorobenzene, o- | 95-50-1 | C6H4Cl2 | 0.5 | 200 | 3 | 50 |
| Dichlorodifluoromethane | 75-71-8 | CCl2F2 | ZR | | | |
| Dichloroethane, 1,2- | 107-06-2 | C2H4Cl2 | ZR | | | |
| Dichloroethane, 1,1- | 75-34-3 | C2H4Cl2 | ZR | | | |
| Dichloroethene, 1,1- | 75-35-4 | C2H2Cl2 | 1.0 | 105 | 5 | 100 |
| Dichloroethene, cis-1,2- | 156-59-2 | C2H2Cl2 | 0.8 | 125 | 4 | 80 |
| Dichloroethene, trans-1,2- | 540-59-0 | C2H2Cl2 | 0.7 | 143 | 4 | 70 |
| Dichloroethylene, 1,2- | 540-59-0 | C2H2Cl2 | 0.8 | 133 | 4 | 75 |
| Dichlorofluoromethane | 75-43-4 | CHFCl2 | ZR | | | |
| Dichloromethane | 75-09-2 | CH2Cl2 | 39 | 3 | 200 | 3900 |
| Dichloropropane, 1,2- | 78-87-5 | C3H6Cl2 | ZR | | | |
| Dichlorotetrafluoroethane, 1,1- | 374-07-2 | C2Cl2F4 | ZR | | | |
| Dichlorotetrafluoroethane, 1,2- | 76-14-2 | C2Cl2F4 | ZR | | | |
| Dicyclopentadiene | 77-73-6 | C10H12 | 0.9 | 110 | 5 | 90 |
| Diesel Fuel | 68334-30-5 | | 0.8 | 130 | 4 | 75 |
| Diethyl ether | 60-29-7 | C4H10O | 0.9 | 110 | 4 | 90 |
| Diethyl maleate | 141-05-9 | C8H12O4 | 2.0 | 50 | 10 | 200 |
| Diethyl phthalate | 84-66-2 | C12H14O4 | 1.0 | 100 | 5 | 100 |
| Diethyl sulphate | 64-67-5 | C4H10SO4 | 3.0 | 33 | 15 | 300 |
| Diethyl sulphide | 352-93-2 | C4H10S | 0.6 | 180 | 3 | 50 |
| Diethylamine | 109-89-7 | C4H11N | 1.0 | 100 | 5 | 100 |
| Diethylaminoethanol, 2- | 100-37-8 | C6H15ON | 2.7 | 40 | 15 | 270 |
| Diethylaminopropylamine, 3- | 104-78-9 | C7H18N2 | 1.0 | 100 | 5 | 100 |
| Difluoroethane, 1,1- | 75-37-6 | C2H4F2 | ZR | | | |
| Difluoroethane, 1,2- | 624-72-6 | C2H4F2 | ZR | | | |
| Difluoromethane | 75-10-5 | CH2F2 | ZR | | | |
| Dihydrogen selenide | 7783-07-5 | H2Se | 1.0 | 100 | 5 | 100 |
| Dihydroxybenzene, 1,2 | 120-80-9 | C6H6O2 | 1.0 | 100 | 5 | 100 |
| Dihydroxybenzene, 1,3 | 108-46-3 | C6H6O2 | 1.0 | 100 | 5 | 100 |
| Diisobutylene | 107-39-1 | C8H16 | 0.6 | 156 | 3 | 60 |
| Diisopropyl ether | 108-20-3 | C6H14O | 0.7 | 150 | 3 | 70 |
| Diisopropylamine | 108-18-9 | C6H15N | 0.7 | 140 | 4 | 70 |
| Diketene | 674-82-8 | C4H4O2 | 2.2 | 45 | 11 | 220 |

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| Gas/ VOC | CAS No. | Formula | Relative response | Relative sensitivity (%) | Typical MDL Mini PID HI (ppb) | Typical MDL Mini PID LO (ppb) |
|-----------------------------------|------------|-----------|-------------------|--------------------------|-------------------------------|-------------------------------|
| Dimethoxymethane | 109-87-5 | C3H8O2 | 1.4 | 71 | 7 | 140 |
| Dimethyl cyclohexane, 1,2- | 583-57-3 | C8H16 | 1.1 | 95 | 5 | 105 |
| Dimethyl disulphide | 624-92-0 | C2H6S2 | 0.2 | 435 | 1 | 23 |
| Dimethyl ether | 115-10-6 | C2H6O | 1.3 | 80 | 7 | 130 |
| Dimethyl phthalate | 131-11-3 | C10H10O4 | 1.0 | 100 | 5 | 100 |
| Dimethyl sulphate | 77-78-1 | C2H6O4S | ZR | | | |
| Dimethyl sulphide | 75-18-3 | C2H6S | 0.5 | 200 | 3 | 50 |
| Dimethylacetamide N,N- | 127-19-5 | C4H9NO | 1.3 | 75 | 7 | 130 |
| Dimethylamine | 124-40-3 | C2H7N | 1.4 | 70 | 7 | 140 |
| Dimethylaminoethanol | 108-01-0 | C4H11NO | 1.5 | 70 | 8 | 150 |
| Dimethylaniline, NN- | 121-69-7 | C8H11N | 0.6 | 167 | 3 | 60 |
| Dimethylbutyl acetate | 108-84-9 | C8H16O2 | 1.6 | 60 | 8 | 160 |
| Dimethylethylamine, NN- | 598-56-1 | C4H11N | 0.8 | 125 | 4 | 80 |
| Dimethylformamide | 68-12-2 | C3H7NO | 0.9 | 110 | 5 | 90 |
| Dimethylheptan-4-one, 2,6- | 108-83-8 | C9H18O | 0.8 | 125 | 4 | 80 |
| Dimethylhydrazine, 1,1- | 57-14-7 | C2H8N2 | 1.0 | 100 | 5 | 100 |
| Dinitrobenzene, m- | 99-65-0 | C6H4N2O4 | 3.0 | 33 | 15 | 300 |
| Dinitrobenzene, o- | 528-29-0 | C6H4N2O4 | ZR | | | |
| Dinitrobenzene, p- | 100-25-4 | C6H4N2O4 | 5.0 | 20 | 25 | 500 |
| Dinonyl phthalate | 84-76-4 | C26H42O4 | 1.0 | 100 | 5 | 100 |
| Dioxane 1,2- | | C4H8O2 | 1.5 | 67 | 8 | 150 |
| Dioxane 1,4- | 123-91-1 | C4H8O2 | 1.5 | 67 | 8 | 150 |
| Dipentene | 138-86-3 | C10H16 | 0.9 | 110 | 5 | 90 |
| Diphenyl ether | 101-84-8 | C12H10O | 0.8 | 125 | 4 | 80 |
| Disulphur decafluoride | 5714-22-7 | S2F10 | ZR | | | |
| Disulphur dichloride | 10025-67-9 | S2Cl2 | 3.0 | 33 | 15 | 300 |
| Di-tert-butyl-p-cresol | 2409-55-4 | C11H16O | 1.0 | 100 | 5 | 100 |
| Divinylbenzene | 1321-74-0 | C10H10 | 0.4 | 250 | 2 | 40 |
| Dodecanol | 112-53-8 | C12H26O | 0.9 | 110 | 5 | 90 |
| Enflurane | 13838-16-9 | C4H2F5ClO | ZR | | | |
| Epichlorohydrin | 106-89-8 | C3H5ClO | 8.0 | 15 | 40 | 800 |
| Epoxypropyl isopropyl ether, 2,3- | 4016-14-2 | C6H12O2 | 1.1 | 90 | 5 | 110 |
| Ethane | 74-84-0 | C2H6 | ZR | | | |
| Ethanol | 64-17-5 | C2H6O | 8.7 | 10 | 45 | 870 |
| Ethanolamine | 141-43-5 | C2H7NO | 3.0 | 33 | 15 | 300 |
| Ethoxy-2-propanol, 1- | 1569-02-4 | C5H10O2 | 2.0 | 50 | 10 | 200 |
| Ethoxyethanol, 2- | 110-80-5 | C4H10O2 | 29.8 | 3 | 150 | 3000 |
| Ethoxyethyl acetate, 2- | 111-15-9 | C6H12O3 | 3.0 | 33 | 15 | 300 |
| Ethyl (S)-(-)-lactate | 97-64-3 | C5H10O3 | 3.0 | 33 | 15 | 300 |
| Ethyl acetate | 141-78-6 | C4H8O2 | 3.6 | 28 | 20 | 360 |
| Ethyl acrylate | 140-88-5 | C5H8O2 | 2.0 | 50 | 10 | 200 |
| Ethyl amine | 75-04-7 | C2H7N | 1.0 | 100 | 5 | 100 |
| Ethyl benzene | 100-41-4 | C8H10 | 0.5 | 185 | 3 | 50 |
| Ethyl butyrate | 105-54-4 | C6H12O2 | 1.0 | 105 | 5 | 100 |
| Ethyl chloroformate | 541-41-3 | C3H5O2Cl | 80 | 1 | 400 | 8300 |
| Ethyl cyanoacrylate | 7085-85-0 | C6H7O2N | 1.5 | 67 | 8 | 150 |
| Ethyl decanoate | 110-38-3 | C12H24O2 | 1.8 | 56 | 10 | 180 |
| Ethyl formate | 109-94-4 | C3H6O2 | 30 | 3 | 150 | 3000 |
| Ethyl hexanoate | 123-66-0 | C8H16O2 | 2.6 | 38 | 15 | 260 |
| Ethyl hexanol, 2- | 105-76-7 | C8H18O | 1.5 | 67 | 8 | 150 |
| Ethyl hexyl acrylate, 2- | 103-11-7 | C11H20O2 | 1.0 | 100 | 5 | 100 |
| Ethyl mercaptan | 75-08-1 | C2H6S | 0.7 | 145 | 3 | 70 |
| Ethyl octanoate | 106-32-1 | C10H20O2 | 2.3 | 40 | 12 | 230 |

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| Gas/ VOC | CAS No. | Formula | Relative response | Relative sensitivity (%) | Typical MDL Mini PID HI (ppb) | Typical MDL Mini PID LO (ppb) |
|-------------------------------------|------------|-----------|-------------------|--------------------------|-------------------------------|-------------------------------|
| Ethylene | 74-85-1 | C2H4 | 8.0 | 13 | 40 | 800 |
| Ethylene dinitrate | 628-96-6 | C2H4O6N2 | ZR | | | |
| Ethylene glycol | 107-21-1 | C2H6O2 | 20.0 | 5 | 100 | 2000 |
| Ethylene oxide | 75-21-8 | C2H4O | 15.0 | 7 | 75 | 1500 |
| Ferrocene | 102-54-5 | C10H10Fe | 0.8 | 125 | 4 | 80 |
| Fluorine | 7782-41-4 | F2 | ZR | | | |
| Fluoroethane | 353-33-6 | C2H5F | ZR | | | |
| Fluoromethane | 593-53-3 | CH3F | ZR | | | |
| Formaldehyde | 50-00-0 | CH2O | ZR | | | |
| Formamide | 75-12-7 | CH3ON | 2.0 | 50 | 10 | 200 |
| Formic acid | 64-18-6 | CH2O2 | ZR | | | |
| Furfural | 98-01-1 | C5H4O2 | 1.4 | 70 | 7 | 140 |
| Furfuryl alcohol | 98-00-0 | C5H6O2 | 2.0 | 50 | 10 | 200 |
| Gasoline vapors | 8006-61-9 | | 1.1 | 95 | 5 | 105 |
| Gasoline vapors | 8006-61-9 | | 0.8 | 125 | 4 | 80 |
| Gasoline vapors 92 octane | 8006-61-9 | | 0.8 | 125 | 4 | 80 |
| Germane | 7782-65-2 | GeH4 | 10.0 | 10 | 50 | 1000 |
| Glutaraldehyde | 111-30-8 | C5H8O2 | 0.9 | 111 | 5 | 90 |
| Halothane | 151-67-7 | CF3CHBrCl | ZR | | | |
| Helium | | He | ZR | | | |
| Heptan-2-one | 110-43-0 | C7H14O | 0.7 | 140 | 4 | 70 |
| Heptan-3-one | 106-35-4 | C7H14O | 0.8 | 133 | 4 | 75 |
| Heptane n- | 142-82-5 | C7H16 | 2.1 | 50 | 10 | 200 |
| Hexachloroethane | 67-72-1 | C2Cl6 | ZR | | | |
| Hexafluoroethane | 76-16-4 | C2F6 | ZR | | | |
| Hexamethyldisilazane, 1,1,1,3,3,3-. | 999-97-3 | C6H19NSi2 | 1.0 | 100 | 5 | 100 |
| Hexamethylcyclotriphospha siloxane. | 107-46-0 | C6H18OSi2 | 0.3 | 350 | 1 | 30 |
| Hexan-2-one | 591-78-6 | C6H12O | 0.8 | 125 | 4 | 80 |
| Hexane n- | 110-54-3 | C6H14 | 4.2 | 25 | 20 | 420 |
| Hexene, 1- | 592-41-6 | C6H12 | 0.9 | 110 | 5 | 90 |
| Hydrazine | 302-01-2 | H4N2 | 3.0 | 33 | 15 | 300 |
| Hydrazoic acid | 7782-79-8 | HN3 | ZR | | | |
| Hydrogen | 1333-74-0 | H2 | ZR | | | |
| Hydrogen bromide | 10035-10-6 | HBr | ZR | | | |
| Hydrogen chloride | 7647-01-0 | HCl | ZR | | | |
| Hydrogen cyanide | 74-90-8 | HCN | ZR | | | |
| Hydrogen fluoride | 7664-39-3 | HF | ZR | | | |
| Hydrogen peroxide | 7722-84-1 | H2O2 | 4.0 | 25 | 20 | 400 |
| Hydrogen sulfide | 7783-06-4 | H2S | 4.0 | 25 | 20 | 400 |
| Hydroquinone | 123-31-9 | C6H6O2 | 0.8 | 125 | 4 | 80 |
| Hydroxypropyl acrylate 2- | 999-61-1 | C6H10O3 | 1.5 | 67 | 8 | 150 |
| Iminodi(ethylamine) 2,2- | 111-40-0 | C4H13N3 | 0.9 | 110 | 5 | 90 |
| Iminodiethanol 2,2'- | 111-42-2 | C4H11NO2 | 1.6 | 60 | 8 | 160 |
| Indene | 95-13-6 | C9H8 | 0.5 | 220 | 2 | 50 |
| Iodine | 7553-56-2 | I2 | 0.2 | 667 | 1 | 15 |
| Iodoform | 75-47-8 | CHI3 | 1.5 | 67 | 8 | 150 |
| Iodomethane | 74-88-4 | CH3I | 0.4 | 250 | 2 | 40 |
| Isoamyl acetate | 123-92-2 | C7H14O2 | 1.6 | 60 | 8 | 160 |
| Isobutane | 75-28-5 | C4H10 | 8.0 | 15 | 40 | 800 |
| Isobutanol | 78-83-1 | C4H10O | 3.5 | 30 | 20 | 350 |
| Isobutyl acetate | 110-19-0 | C6H12O2 | 2.3 | 45 | 10 | 230 |
| Isobutyl acrylate | 106-63-8 | C7H12O2 | 1.3 | 80 | 7 | 130 |
| Isobutylene | 115-11-7 | C4H8 | 1.0 | 100 | 5 | 100 |

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| Gas/ VOC | CAS No. | Formula | Relative response | Relative sensitivity (%) | Typical MDL Mini PID HI (ppb) | Typical MDL Mini PID LO (ppb) |
|---------------------------------|------------|-----------|-------------------|--------------------------|-------------------------------|-------------------------------|
| Isobutyraldehyde | 78-84-2 | C4H8O | 1.2 | 80 | 6 | 120 |
| Isocyanates, all | | | NV | | | |
| Isodecanol | 25339-17-7 | C10H22O | 0.9 | 110 | 5 | 90 |
| Isoflurane | 26675-46-7 | C3H2ClF5O | ZR | | | |
| Isononanol | 2452-97-9 | C9H20O | 1.5 | 67 | 8 | 150 |
| Isooctane | 565-75-3 | C8H18 | 1.1 | 90 | 5 | 100 |
| Isooctanol | 26952-21-6 | C8H18O | 1.7 | 60 | 9 | 170 |
| Isopentane | 78-78-4 | C5H12 | 6.0 | 20 | 30 | 600 |
| Isophorone | 78-59-1 | C9H14O | 0.8 | 133 | 4 | 75 |
| Isoprene | 78-79-5 | C5H8 | 0.7 | 140 | 3 | 70 |
| Isopropanol | 67-63-0 | C3H8O | 4.4 | 20 | 22 | 440 |
| Isopropyl acetate | 108-21-4 | C5H10O2 | 2.2 | 50 | 10 | 220 |
| Isopropyl chloroformate | 108-23-6 | C4H7O2Cl | 1.6 | 60 | 8 | 160 |
| Jet Fuel JP-4 | | | 0.8 | 133 | 4 | 75 |
| Jet Fuel JP-5 | | | 0.7 | 150 | 3 | 60 |
| Jet Fuel JP-8 | | | 0.7 | 150 | 3 | 60 |
| Kerosene | 8008-20-6 | | 0.8 | 120 | 4 | 90 |
| Ketene | 463-51-4 | C2H2O | 3.0 | 33 | 15 | 300 |
| Liquefied petroleum gas | 68476-85-7 | | ZR | | | |
| Maleic anhydride | 108-31-6 | C4H2O3 | 2.0 | 50 | 10 | 200 |
| Mercaptoacetic acid | 68-11-1 | C2H4O2S | 1.0 | 100 | 5 | 100 |
| Mercury | 7439-97-6 | Hg | NV | | | |
| Mercury alkyls | | | NV | | | |
| Mesitylene | 108-67-8 | C9H12 | 0.3 | 300 | 2 | 30 |
| Methacrylic acid | 79-41-4 | C4H6O2 | 2.3 | 40 | 12 | 230 |
| Methacrylonitrile | 126-98-7 | C4H5N | 5.0 | 20 | 25 | 500 |
| Methane | 74-82-8 | CH4 | ZR | | | |
| Methanol | 67-56-1 | CH4O | 200 | 1 | 1000 | 20000 |
| Methoxyethanol, 2- | 109-86-4 | C3H8O2 | 2.7 | 40 | 15 | 270 |
| Methoxyethoxyethanol, 2- | 111-77-3 | C5H12O3 | 1.4 | 70 | 7 | 140 |
| Methoxymethylmethoxy-2-propanol | 34590-94-8 | C7H16O3 | 1.3 | 80 | 7 | 130 |
| Methoxypropan-2-ol | 107-98-2 | C4H10O2 | 3.0 | 33 | 15 | 300 |
| Methoxypropyl acetate | 108-65-6 | C6H12O3 | 1.2 | 80 | 6 | 120 |
| Methyl acetate | 79-20-9 | C3H6O2 | 5.2 | 20 | 25 | 500 |
| Methyl acrylate | 96-33-3 | C4H6O2 | 3.4 | 30 | 17 | 340 |
| Methyl bromide | 74-83-9 | CH3Br | 1.9 | 50 | 10 | 190 |
| Methyl cyanoacrylate | 137-05-3 | C5H5O2N | 5.0 | 20 | 25 | 500 |
| Methyl ethyl ketone | 78-93-3 | C4H8O | 0.8 | 130 | 4 | 80 |
| Methyl ethyl ketone peroxides | 1338-23-4 | C8H18O2 | 0.8 | 125 | 4 | 80 |
| Methyl formate | 107-31-3 | C2H4O2 | ZR | | | |
| Methyl isobutyl ketone | 108-10-1 | C6H12O | 0.8 | 125 | 4 | 80 |
| Methyl isocyanate | 624-83-9 | C2H3NO | ZR | | | |
| Methyl isothiocyanate | 556-61-6 | C2H3NS | 0.6 | 167 | 3 | 60 |
| Methyl mercaptan | 74-93-1 | CH4S | 0.7 | 140 | 4 | 70 |
| Methyl methacrylate | 80-62-6 | C5H8O2 | 1.6 | 60 | 8 | 160 |
| Methyl propyl ketone | 107-87-9 | C5H10O | 0.8 | 130 | 4 | 80 |
| Methyl salicylate | 119-36-8 | C8H8O3 | 1.2 | 80 | 6 | 120 |
| Methyl sulphide | 75-18-3 | C2H6S | 0.5 | 200 | 3 | 50 |
| Methyl t-butyl ether | 1634-04-4 | C5H12O | 0.8 | 125 | 4 | 80 |
| Methyl-2-propen-1-ol, 2- | 51-42-8 | C4H8O | 1.1 | 90 | 5 | 100 |
| Methyl-2-pyrrolidinone, N- | 872-50-4 | C5H9NO | 0.9 | 110 | 5 | 90 |
| Methyl-4,6-dinitrophenol, 2- | 534-52-1 | C7H6N2O5 | 3.0 | 33 | 15 | 300 |
| Methyl-5-hepten-2-one, 6- | 110-93-0 | C8H14O | 0.8 | 125 | 4 | 80 |

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|---------------------------------------|------------|----------|-------------------|--------------------------|-------------------------------|-------------------------------|
| Methylamine | 74-89-5 | CH5N | 1.4 | 70 | 7 | 140 |
| Methylbutan-1-ol, 3- | 123-51-3 | C5H12O | 3.4 | 30 | 17 | 340 |
| Methylcyclohexane | 108-87-2 | C7H14 | 1.1 | 90 | 6 | 110 |
| Methylcyclohexanol, 4- | 589-91-3 | C7H14O | 2.4 | 40 | 12 | 240 |
| Methylcyclohexanone 2- | 583-60-8 | C7H12O | 1.0 | 100 | 5 | 100 |
| Methylheptan-3-one, 5- | 541-85-5 | C8H16O | 0.8 | 133 | 4 | 75 |
| Methylhexan-2-one, 5- | 110-12-3 | C7H14O | 0.8 | 133 | 4 | 75 |
| Methylhydrazine | 60-34-4 | CH6N2 | 1.3 | 80 | 7 | 130 |
| Methyl-N-2,4, 6-tetranitroaniline, N- | 479-45-8 | C7H5N5O8 | 3.0 | 33 | 15 | 300 |
| Methylpent-3-en-2-one, 4- | 141-79-7 | C6H10O | 0.7 | 140 | 4 | 70 |
| Methylpentan-2-ol, 4- | 108-11-2 | C6H14O | 2.8 | 40 | 14 | 280 |
| Methylpentane-2,4-diol, 2- | 107-41-5 | C6H14O2 | 4.0 | 25 | 20 | 400 |
| Methylpropan-2-ol, 2- | 75-65-0 | C4H10O | 3.5 | 30 | 18 | 350 |
| Methylstyrene | 25013-15-4 | C9H10 | 0.5 | 200 | 3 | 50 |
| Mineral oil | 8042-47-5 | | 0.8 | 125 | 4 | 80 |
| Mineral spirits | 64475-85-0 | | 0.8 | 125 | 4 | 80 |
| Naphthalene | 91-20-3 | C10H8 | 0.4 | 230 | 2 | 45 |
| Nitric oxide | 10102-43-9 | NO | 8.0 | 15 | 40 | 800 |
| Nitroaniline 4- | 100-01-6 | C6H6N2O2 | 0.8 | 125 | 4 | 80 |
| Nitrobenzene | 98-95-3 | C6H5NO2 | 1.7 | 60 | 10 | 170 |
| Nitroethane | 79-24-3 | C2H5NO2 | ZR | | | |
| Nitrogen dioxide | 10102-44-0 | NO2 | 10.0 | 10 | 50 | 1000 |
| Nitrogen trichloride | 10025-85-1 | NC13 | 1.0 | 100 | 5 | 100 |
| Nitrogen trifluoride | 7783-54-2 | NF3 | ZR | | | |
| Nitromethane | 75-52-5 | CH3NO2 | ZR | | | |
| Nitropropane, 1- | 108-03-2 | C3H7NO2 | ZR | | | |
| Nitropropane, 2- | 79-46-9 | C3H7NO2 | ZR | | | |
| Nitrous oxide | 10024-97-2 | N2O | ZR | | | |
| Nonane, n- | 111-84-2 | C9H20 | 1.3 | 80 | 6 | 130 |
| Norbornadiene, 2,5- | 121-46-0 | C7H8 | 0.6 | 167 | 3 | 60 |
| Octachloronaphthalene | 2234-13-1 | C10Cl8 | 1.0 | 100 | 5 | 100 |
| Octane, n- | 111-65-9 | C8H18 | 1.6 | 60 | 8 | 160 |
| Octene, 1- | 111-66-0 | C8H16 | 0.7 | 140 | 3 | 70 |
| Oxalic acid | 144-62-7 | C2H2O4 | ZR | | | |
| Oxalonitrile | 460-19-5 | C2N2 | ZR | | | |
| Oxydiethanol 2,2- | 111-46-6 | C4H10O3 | 4.0 | 25 | 20 | 400 |
| Oxygen | | O2 | ZR | | | |
| Ozone | 10028-15-6 | O3 | ZR | | | |
| Paraffin wax, fume | 8002-74-2 | | 1.0 | 100 | 5 | 100 |
| Paraffins, normal | 64771-72-8 | | 1.0 | 105 | 5 | 100 |
| Pentacarbonyl iron | 13463-40-6 | FeC5O5 | 1.0 | 100 | 5 | 100 |
| Pentachloroethane | 76-01-7 | C2HCl5 | ZR | | | |
| Pentachlorofluoroethane | 354-56-3 | C2Cl5F | ZR | | | |
| Pentafluoroethane | 354-33-6 | C2HF5 | ZR | | | |
| Pentan-2-one | 107-87-9 | C5H10O | 0.8 | 125 | 4 | 80 |
| Pentan-3-one | 96-22-0 | C5H10O | 0.8 | 125 | 4 | 80 |
| Pentandione, 2,4- | 123-54-6 | C5H8O2 | 0.8 | 133 | 4 | 75 |
| Pentane, n- | 109-66-0 | C5H12 | 7.9 | 15 | 40 | 800 |
| Peracetic acid | 79-21-0 | C2H4O3 | 2.0 | 50 | 10 | 200 |
| Perchloryl fluoride | 7616-94-6 | ClO3F | ZR | | | |
| Perfluoropropane | 76-19-7 | C3F8 | ZR | | | |
| Petroleum ether | | | 0.9 | 110 | 5 | 90 |
| Phenol | 108-95-2 | C6H6O | 1.2 | 85 | 6 | 120 |

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| Gas/ VOC | CAS No. | Formula | Relative response | Relative sensitivity (%) | Typical MDL Mini PID HI (ppb) | Typical MDL Mini PID LO (ppb) |
|--|------------|-----------|-------------------|--------------------------|-------------------------------|-------------------------------|
| Phenyl propene, 2- | 98-83-9 | C9H10 | 0.4 | 230 | 2 | 45 |
| Phenyl-2,3-epoxypropyl ether | 122-60-1 | C9H10O2 | 0.8 | 125 | 4 | 80 |
| Phenylenediamine, p- | 106-50-3 | C6H8N2 | 0.6 | 167 | 3 | 60 |
| Phosgene | 75-44-5 | COCl2 | ZR | | | |
| Phosphine | 7803-51-2 | PH3 | 2.0 | 50 | 10 | 200 |
| Picoline, 3- | 108-99-6 | C6H7N | 0.9 | 110 | 5 | 90 |
| Pinene, alpha | 80-56-8 | C10H16 | 0.3 | 315 | 2 | 30 |
| Pinene, beta | 127-91-3 | C10H16 | 0.3 | 315 | 2 | 30 |
| Piperidine | 110-89-4 | C5H11N | 0.9 | 110 | 5 | 90 |
| Piperylene | 504-60-9 | C5H8 | 0.7 | 150 | 3 | 67 |
| Prop-2-yn-1-ol | 107-19-7 | C3H4O | 1.3 | 80 | 7 | 130 |
| Propan-1-ol | 71-23-8 | C3H8O | 4.8 | 20 | 25 | 480 |
| Propane | 74-98-6 | C3H8 | ZR | | | |
| Propane-1,2-diol, total | 57-55-6 | C3H8O2 | 10.0 | 10 | 50 | 1000 |
| Propene | 115-07-1 | C3H6 | 1.4 | 70 | 7 | 140 |
| Propionaldehyde | 123-38-6 | C3H6O | 1.7 | 60 | 8 | 169 |
| Propionic acid | 79-09-4 | C3H6O2 | 8.0 | 15 | 40 | 800 |
| Propyl acetate, n- | 109-60-4 | C5H10O2 | 2.5 | 40 | 13 | 250 |
| Propylene dinitrate | 6423-43-4 | C3H6N2O6 | ZR | | | |
| Propylene oxide | 75-56-9 | C3H6O | 7.0 | 15 | 35 | 700 |
| Propyleneimine | 75-55-8 | C3H7N | 1.3 | 80 | 7 | 130 |
| Pyridine | 110-86-1 | C5H5N | 0.8 | 133 | 4 | 75 |
| Pyridylamine 2- | 504-29-0 | C5H6N2 | 0.8 | 125 | 4 | 80 |
| Silane | 7803-62-5 | SiH4 | ZR | | | |
| Sodium fluoroacetate | 62-74-8 | C2H2O2FNa | ZR | | | |
| Styrene | 100-42-5 | C8H8 | 0.4 | 230 | 2 | 50 |
| Sulphur dioxide | 7446-09-5 | SO2 | ZR | | | |
| Sulphur hexafluoride | 2551-62-4 | SF6 | ZR | | | |
| Sulphur tetrafluoride | 7783-60-0 | SF4 | ZR | | | |
| Sulphuric acid | 7664-93-9 | H2SO4 | ZR | | | |
| Sulphuryl fluoride | 2699-79-8 | SO2F2 | ZR | | | |
| Terphenyls | | C18H14 | 0.6 | 167 | 3 | 60 |
| Terpinolene | 586-62-9 | C10H16 | 0.5 | 210 | 2 | 50 |
| Tert-butanol | 75-65-0 | C4H10O | 2.6 | 40 | 15 | 260 |
| Tetrabromoethane, 1,1,2,2- | 79-27-6 | C2H2Br4 | 2.0 | 50 | 10 | 200 |
| Tetracarbonylnickel | 13463-39-3 | NiC4O4 | 1.0 | 100 | 5 | 100 |
| Tetrachloro-1,2-difluoroethane, 1,1,2,2- | 76-12-0 | C2Cl4F2 | ZR | | | |
| Tetrachloro-1-fluoroethane, 1,1,2,2- | 354-14-3 | C2HCl4F | ZR | | | |
| Tetrachloro-2,2-difluoroethane, 1,1,1,2- | 76-11-9 | C2Cl4F2 | ZR | | | |
| Tetrachloro-2-fluoroethane, 1,1,1,2- | 354-11-0 | C2HCl4F | ZR | | | |
| Tetrachloroethane, 1,1,1,2- | 630-20-6 | C2H2Cl4 | ZR | | | |
| Tetrachloroethane, 1,1,2,2- | 79-34-5 | C2H2Cl4 | ZR | | | |
| Tetrachloroethylene | 127-18-4 | C2Cl4 | 0.7 | 140 | 4 | 70 |
| Tetrachloronaphthalenes, all isomers | 20020-02-4 | C10H4Cl4 | 1.0 | 100 | 5 | 100 |
| Tetraethyl orthosilicate | 78-10-4 | C8H20O4Si | 2.0 | 50 | 10 | 200 |
| Tetraethyllead | 78-00-2 | C8H20Pb | ZR | | | |
| Tetrafluoroethane, 1,1,1,2- | 811-97-2 | C2H2F4 | ZR | | | |
| Tetrafluoroethane, 1,1,2,2- | 359-35-3 | C2H2F4 | ZR | | | |
| Tetrafluoroethylene | 116-14-3 | C2F4 | 1.0 | 100 | 5 | 100 |
| Tetrafluromethane | 75-73-0 | CF4 | ZR | | | |

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| Gas/ VOC | CAS No. | Formula | Relative response | Relative sensitivity (%) | Typical MDL Mini PID HI (ppb) | Typical MDL Mini PID LO (ppb) |
|--------------------------------------|-----------|-----------|-------------------|--------------------------|-------------------------------|-------------------------------|
| Tetrahydrofuran | 109-99-9 | C4H8O | 1.6 | 65 | 8 | 150 |
| Tetramethyl orthosilicate | 681-84-5 | C4H12O4Si | ZR | | | |
| Tetramethyl succinonitrile | 3333-52-6 | C8H12N2 | 1.0 | 100 | 5 | 100 |
| Therminol | | | 1.0 | 100 | 5 | 100 |
| Thionyl chloride | 7719-09-7 | SOCl2 | ZR | | | |
| Toluene | 108-88-3 | C7H8 | 0.5 | 200 | 3 | 50 |
| Toluene-2,4-diisocyanate | 584-84-9 | C9H6N2O2 | 1.6 | 60 | 8 | 160 |
| Toluenesulphonyl chloride, p- | 98-59-9 | C7H7SO2Cl | 3.0 | 33 | 15 | 300 |
| Toluidine, o- | 95-53-4 | C7H9N | 0.5 | 200 | 3 | 50 |
| Tributyl phosphate | 126-73-8 | C12H27O4P | 5.0 | 20 | 25 | 500 |
| Tributylamine | 102-82-9 | C12H27N | 1.0 | 100 | 5 | 100 |
| Trichloro-1,1-difluoroethane, 1,2,2- | 354-21-2 | C2HCl3F2 | ZR | | | |
| Trichloro-1,2-difluoroethane, 1,1,2- | 354-15-4 | C2HCl3F2 | ZR | | | |
| Trichloro-2,2-difluoroethane, 1,1,1- | 354-12-1 | C2HCl3F2 | ZR | | | |
| Trichloro-2-fluoroethane, 1,1,2- | 359-28-4 | C2H2Cl3F | ZR | | | |
| Trichlorobenzene 1,2,4- | 120-82-1 | C6H3Cl3 | 0.6 | 180 | 3 | 50 |
| Trichloroethane, 1,1,1- | 71-55-6 | C2H3Cl3 | ZR | | | |
| Trichloroethane, 1,1,2- | 79-00-5 | C2H3Cl3 | ZR | | | |
| Trichloroethylene | 79-01-6 | C2HCl3 | 0.7 | 150 | 3 | 65 |
| Trichlorofluoromethane | 75-69-4 | CCl3F | ZR | | | |
| Trichloronitromethane | 76-06-2 | CCl3NO2 | ZR | | | |
| Trichlorophenoxyacetic acid, 2,4,5- | 93-76-5 | C8H5O3Cl3 | 1.0 | 100 | 5 | 100 |
| Trichloropropane 1,2,3- | 96-18-4 | C3H5Cl3 | ZR | | | |
| Trichlorotrifluoroethane, 1,1,1- | 354-58-5 | C2Cl3F3 | ZR | | | |
| Trichlorotrifluoroethane, 1,1,2- | 76-13-1 | C2Cl3F3 | ZR | | | |
| Triethylamine | 121-44-8 | C6H15N | 0.9 | 110 | 5 | 90 |
| Trifluoroethane, 1,1,1- | 420-46-2 | C2H3F3 | ZR | | | |
| Trifluoroethane, 1,1,2- | 430-66-0 | C2H3F3 | ZR | | | |
| Trifluoroethanol, 2,2,2- | 75-89-8 | C2H3F3O | ZR | | | |
| Trifluoromethane | 75-46-7 | CHF3 | ZR | | | |
| Trimethylamine | 53-50-3 | C3H9N | 0.5 | 200 | 3 | 50 |
| Trimethylbenzene mixtures | | C9H12 | 0.3 | 300 | 2 | 35 |
| Trimethylbenzene, 1,3,5- | 108-67-8 | C9H12 | 0.3 | 300 | 2 | 35 |
| Trinitrotoluene 2,4,6- | 118-96-7 | C7H5N3O6 | ZR | | | |
| Turpentine | 8006-64-2 | C10H16 | 0.6 | 167 | 3 | 60 |
| TVOC | | | 1.0 | 100 | 5 | 100 |
| Undecane, n- | 1120-21-4 | C11H24 | 0.9 | 110 | 5 | 100 |
| Vinyl acetate | 108-05-2 | C4H6O2 | 1.1 | 90 | 6 | 110 |
| Vinyl bromide | 593-60-2 | C2H3Br | 1.0 | 100 | 5 | 100 |
| Vinyl chloride | 75-01-4 | C2H3Cl | 2.1 | 50 | 10 | 200 |
| Vinyl-2-pyrrolidinone, 1- | 88-12-0 | C6H9NO | 0.9 | 110 | 5 | 90 |
| Xylene mixed isomers | 1330-20-7 | C8H10 | 0.4 | 230 | 2 | 40 |
| Xylene, m- | 108-38-3 | C8H10 | 0.4 | 230 | 2 | 50 |
| Xylene, o- | 95-47-6 | C8H10 | 0.6 | 167 | 3 | 60 |
| Xylene, p- | 106-42-3 | C8H10 | 0.6 | 180 | 3 | 50 |
| Xyliidine, all | 1300-73-8 | C8H11N | 0.7 | 140 | 4 | 70 |

Advanced Gas Sensing Technologies

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