

## CORRECTION FACTORS, IONIZATION POTENTIALS, AND CALIBRATION CHARACTERISTICS

### Correction Factors and Ionization Potentials

RAE Systems PIDs can be used for the detection of a wide variety of gases that exhibit different responses. In general, any compound with ionization potential (IP) lower than that of the lamp photons can be measured.\* The best way to calibrate a PID to different compounds is to use a standard of the gas of interest. However, correction factors have been determined that enable the user to quantify a large number of chemicals using only a single calibration gas, typically isobutylene. In our PIDs, correction factors can be used in one of three ways:

- 1) Calibrate the unit with isobutylene in the usual fashion to read in isobutylene equivalents. Manually multiply the reading by the correction factor (CF) to obtain the concentration of the gas being measured.
- 2) Calibrate the unit with isobutylene in the usual fashion to read in isobutylene equivalents. Call up the correction factor from the instrument memory or download it from a personal computer and then call it up. The unit will then read directly in units of the gas of interest.
- 3) Calibrate the unit with isobutylene, but input an equivalent, "corrected" span gas concentration when prompted for this value. The unit will then read directly in units of the gas of interest.

\* Some inorganic compounds like  $H_2O_2$  and  $NO_2$  give weak response at photon energies well above those of their ionization potentials.

### Example 1:

With the unit calibrated to read isobutylene equivalents, the reading is 10 ppm with a 10.6 eV lamp. The gas being measured is butyl acetate, which has a correction factor of 2.6. Multiplying 10 by 2.6 gives an adjusted butyl acetate value of 26 ppm. Similarly, if the gas being measured were trichloroethylene (CF = 0.54), the adjusted value with a 10 ppm reading would be 5.4 ppm.

### Example 2:

With the unit calibrated to read isobutylene equivalents, the reading is 100 ppm with a 10.6 eV lamp. The gas measured is m-xylene (CF = 0.43). After downloading this factor, the unit should read about 43 ppm when exposed to the same gas, and thus read directly in m-xylene values.

### Example 3:

The desired gas to measure is ethylene dichloride (EDC). The CF is 0.6 with an 11.7 eV lamp. During calibration with 100 ppm isobutylene, insert 0.6 times 100, or 60 at the prompt for the calibration gas concentration. The unit then reads directly in EDC values.

### Conversion to $mg/m^3$

To convert from ppm to  $mg/m^3$ , use the following formula:

$$\text{Conc. (mg/m}^3\text{)} = \frac{[\text{Conc. (ppmv)} \times \text{mol. wt. (g/mole)}]}{\text{molar gas volume (L)}}$$

For air at 25 °C (77 °F), the molar gas volume is 24.4 L/mole and the formula reduces to:

$$\text{Conc. (mg/m}^3\text{)} = \text{Conc. (ppmv)} \times \text{mol. wt. (g/mole)} \times 0.041$$

For example, if the instrument is calibrated with a gas standard in ppmv, such as 100 ppm isobutylene, and the user wants to display to read in  $mg/m^3$  of hexane, whose m.w. is 86 and CF is 4.3, the overall correction factor would be  $4.3 \times 86 \times 0.041$  equals 15.2.

### Correction Factors for Mixtures

The correction factor for a mixture is calculated from the sum of the mole fractions  $X_i$  of each component divided by their respective correction factors  $CF_i$ :

$$CF_{\text{mix}} = 1 / (X_1/CF_1 + X_2/CF_2 + X_3/CF_3 + \dots X_i/CF_i)$$

Thus, for example, a vapor phase mixture of 5% benzene and 95% n-hexane would have a  $CF_{\text{mix}}$  of  $CF_{\text{mix}} = 1 / (0.05/0.53 + 0.95/4.3) = 3.2$ . A reading of 100 would then correspond to 320 ppm of the total mixture, comprised of 16 ppm benzene and 304 ppm hexane.

For a spreadsheet to compute the correction factor and TLV of a mixture see the appendix at the end of the CF table.

### TLVs and Alarm Limits for Mixtures

The correction factor for mixtures can be used to set alarm limits for mixtures. To do this one first needs to calculate the exposure limit for the mixture. The Threshold Limit Value (TLV) often defines exposure limits. The TLV for the mixture is calculated in a manner similar to the CF calculation:

$$TLV_{\text{mix}} = 1 / (X_1/TLV_1 + X_2/TLV_2 + X_3/TLV_3 + \dots X_i/TLV_i)$$

In the above example, the 8-h TLV for benzene is 0.5 ppm and for n-hexane 50 ppm. Therefore the TLV of the mixture is  $TLV_{\text{mix}} = 1 / (0.05/0.5 + 0.95/50) = 8.4$  ppm, corresponding to 8.0 ppm hexane and 0.4 ppm benzene. For an instrument calibrated on isobutylene, the reading corresponding to the TLV is:

$$\text{Alarm Reading} = TLV_{\text{mix}} / CF_{\text{mix}} = 8.4 / 3.2 = 2.6 \text{ ppm}$$

A common practice is to set the lower alarm limit to half the TLV, and the higher limit to the TLV. Thus, one would set the alarms to 1.3 and 2.6 ppm, respectively.

## Calibration Characteristics

a) **Flow Configuration.** PID response is essentially independent of gas flow rate as long as it is sufficient to satisfy the pump demand. Four main flow configurations are used for calibrating a PID:

- 1) **A pressurized gas cylinder (Fixed-flow regulator):** The flow rate of the regulator should match the flow demand of the instrument pump or be slightly higher.
- 2) **A pressurized gas cylinder (Demand-flow regulator):** A demand-flow regulator better matches pump speed differences, but results in a slight vacuum during calibration and thus slightly high readings.
- 3) **A collapsible gas bag:** The instrument will draw the calibration gas from the bag at its normal flow rate, as long as the bag valve is large enough. The bag should be filled with enough gas to allow at least one minute of flow (~ 0.6 L for a MiniRAE, ~0.3 L for MultiRAE).
- 4) **The T (or open tube) method:** The T method uses a T-junction with gas flow higher than the pump draw. The gas supply is connected to one end of the T, the instrument inlet is connected to a second end of the T, and excess gas flow escapes through the third, open end of the T. To prevent ambient air mixing, a long tube should be connected to the open end, or a high excess rate should be used. Alternatively, the instrument probe can be inserted into an open tube slightly wider than the probe. Excess gas flows out around the probe.

The first two cylinder methods are the most efficient in terms of gas usage, while the bag and T methods give slightly more accurate results because they match the pump flow better.

- b) **Pressure.** Pressures deviating from atmospheric pressure affect the readings by altering gas concentration and pump characteristics. It is best to calibrate with the instrument and calibration gas at the same pressure as each other and the sample gas. (Note that the cylinder pressure is not relevant because the regulator reduces the pressure to ambient.) If the instrument is calibrated at atmospheric pressure in one of the flow configurations described above, then 1) pressures slightly above ambient are acceptable but high pressures can damage the pump and 2) samples under vacuum may give low readings if air leaks into the sample train.
- c) **Temperature.** Because temperature effects gas density and concentration, the temperature of the calibration gas and instrument should be as close as possible to the ambient temperature where the unit will be used. We recommend that the temperature of the calibration gas be within the instrument's operating range of 14°-104° F (-10°- 40° C). Also, during actual measurements, the instrument should be kept at the same or higher temperature than the sample temperature to avoid condensation in the unit.
- d) **Matrix.** The matrix gas of the calibration compound and VOC sample is significant. Some common matrix components, such as methane and water vapor can affect the

VOC signal. PIDs are most commonly used for monitoring VOCs in air, in which case the preferred calibration gas matrix is air. For a MiniRAE, methane, methanol, and water vapor reduce the response by about 20% when their concentration is 15,000 ppm and by about 40% at 30,000 ppm. Despite earlier reports of oxygen effects, RAE PID responses with 10.6 eV lamps are independent of oxygen concentration, and calibration gases in a pure nitrogen matrix can be used. H<sub>2</sub> and CO<sub>2</sub> up to 5 volume % also have no effect.

- e) **Concentration.** Although RAE Systems PIDs have electronically linearized output, it is best to calibrate in a concentration range close to the actual measurement range. For example, 100 ppm standard gas for anticipated vapors of 0 - 250 ppm, and 500 ppm standard for expected concentrations of 250 - 1000 ppm.
- f) **Filters.** Filters affect flow and pressure conditions and therefore all filters to be used during sampling should also be in place during calibration. Using a water trap (hydrophobic filter) greatly reduces the chances of drawing water aerosols or dirt particles into the instrument. Regular filter replacements are recommended because dirty filters can adsorb VOCs and cause slower response time and shifts in calibration.

### Table Abbreviations:

- CF** =Correction Factor (multiply by reading to get corrected value for the compound when calibrated to isobutylene)
- NR** =No Response
- IP** =Ionization Potential (values in parentheses are not well established)
- C** =Confirmed Value; all others are preliminary or estimated values and are subject to change
- ne** =Not Established ACGIH 8-hr. TWA
- C##** =Ceiling value, given where 8-hr.TWA is not available

### Disclaimer:

Actual readings may vary with age and cleanliness of lamp, relative humidity, and other factors. For accurate work, the instrument should be calibrated regularly under the operating conditions used. The correction factors in this table were measured in dry air at room temperature.

### Updates:

The values that are indicated by a "plus" sign in the "C" columns are confirmed values; all others are preliminary and subject to change. Watch for updates of this table on the Internet at <http://www.raesystems.com>

*IP data are taken from the CRC Handbook of Chemistry and Physics, 73rd Edition, D.R. Lide (Ed.), CRC Press (1993) and NIST Standard Ref. Database 19A, NIST Positive Ion Energetics, Vers. 2.0, Lias, et.al., U.S. Dept. Commerce (1993). Exposure limits (8-h TWA and Ceiling Values) are from the 1997 ACGIH TLVs and BEIs, Threshold Limit Values for Chemical Substances and Physical Agents and Biological Exposure Indices. ACGIH, Cincinnati, OH 1997.*

Compound Name	Synonym/Abbreviation	CAS No.	Formula	9.8	C	10.6	C	11.7	C	IP (eV)	TWA
Acetaldehyde		75-07-0	C <sub>2</sub> H <sub>4</sub> O	NR	+	6	+	3.3	+	10.23	C25
Acetic Acid	Ethanoic Acid	64-19-7	C <sub>2</sub> H <sub>4</sub> O <sub>2</sub>	NR	+	22	+	2.6	+	10.66	10
Acetic Anhydride	Ethanoic Acid Anhydride	108-24-7	C <sub>4</sub> H <sub>6</sub> O <sub>3</sub>	NR	+	6.1	+	2.0	+	10.14	5
Acetone	2-Propanone	67-64-1	C <sub>3</sub> H <sub>6</sub> O	1.2	+	1.1	+	1.4	+	9.71	500
Acetonitrile	Methyl cyanide, Cyanomethane	75-05-8	C <sub>2</sub> H <sub>3</sub> N					100		12.19	40
Acetylene	Ethyne	74-86-2	C <sub>2</sub> H <sub>2</sub>					2		11.40	ne
Acrolein	Propenal	107-02-8	C <sub>3</sub> H <sub>4</sub> O	42	+	3.9	+	1.4	+	10.10	0.1
Acrylic Acid	Propenoic Acid	79-10-7	C <sub>3</sub> H <sub>4</sub> O <sub>2</sub>			12	+	2.0	+	10.60	2
Acrylonitrile	Propenenitrile	107-13-1	C <sub>3</sub> H <sub>3</sub> N			NR	+	1.2	+	10.91	2
Allyl alcohol		107-18-6	C <sub>3</sub> H <sub>6</sub> O			2.4	+	1.7		9.67	2
Allyl chloride	3-Chloropropene	107-05-1	C <sub>3</sub> H <sub>5</sub> Cl			4.3		0.7		9.9	1
Ammonia		7664-41-7	H <sub>3</sub> N	NR	+	9.7	+	5.7	+	10.16	25
Amyl acetate	mix of n-Pentyl acetate & 2-Methylbutyl acetate	628-63-7	C <sub>7</sub> H <sub>14</sub> O <sub>2</sub>	11	+	2.3	+	0.95	+	<9.9	100
Amyl alcohol	1-Pentanol	75-85-4	C <sub>5</sub> H <sub>12</sub> O			5				10.00	ne
Aniline	Aminobenzene	62-53-3	C <sub>7</sub> H <sub>7</sub> N	0.50	+	0.48	+	0.47	+	7.72	2
Anisole	Methoxybenzene	100-66-3	C <sub>7</sub> H <sub>8</sub> O			0.8				8.21	ne
Arsine	Arsenic trihydride	7784-42-1	AsH <sub>3</sub>			1.9	+			9.89	0.05
Benzaldehyde		100-52-7	C <sub>7</sub> H <sub>6</sub> O					1		9.49	ne
Benzene		71-43-2	C <sub>6</sub> H <sub>6</sub>	0.55	+	0.53	+	0.6	+	9.25	0.5
Benzonitrile	Cyanobenzene	100-47-0	C <sub>7</sub> H <sub>5</sub> N			1.6				9.62	ne
Benzyl alcohol	α-Hydroxytoluene, Hydroxymethylbenzene, Benzenemethanol	100-51-6	C <sub>7</sub> H <sub>8</sub> O	1.4	+	1.1	+	0.9	+	8.26	ne
Benzyl chloride	α-Chlorotoluene, Chloromethylbenzene	100-44-7	C <sub>7</sub> H <sub>7</sub> Cl	0.7	+	0.6	+	0.5	+	9.14	1
Benzyl formate	Formic acid benzyl ester	104-57-4	C <sub>8</sub> H <sub>8</sub> O <sub>2</sub>	0.9	+	0.73	+	0.66	+		ne
Boron trifluoride		7637-07-2	BF <sub>3</sub>	NR		NR		NR		15.5	C1
Bromine		7726-95-6	Br <sub>2</sub>	NR	+	1.30	+	0.74	+	10.51	0.1
Bromobenzene		108-86-1	C <sub>6</sub> H <sub>5</sub> Br			0.6		0.5		8.98	ne
2-Bromoethyl methyl ether		6482-24-2	C <sub>3</sub> H <sub>7</sub> OBr			0.84	+			~10	ne
Bromoform	Tribromomethane	75-25-2	CHBr <sub>3</sub>	NR	+	2.5	+	0.5	+	10.48	0.5
Bromopropane, 1-	n-Propyl bromide	106-94-5	C <sub>3</sub> H <sub>7</sub> Br	150	+	1.5	+	0.6	+	10.18	ne
Butadiene	1,3-Butadiene, Vinyl ethylene	106-99-0	C <sub>4</sub> H <sub>6</sub>	0.8		0.85	+	1.1		9.07	2
Butadiene diepoxide, 1,3-	1,2,3,4-Diepoxybutane	298-18-0	C <sub>4</sub> H <sub>6</sub> O <sub>2</sub>	25	+	3.5	+	1.2		~10	ne
Butane		106-97-8	C <sub>4</sub> H <sub>10</sub>			67	+	1.2		10.53	800
Butanol, 1-	Butyl alcohol, n-Butanol	71-36-3	C <sub>4</sub> H <sub>10</sub> O	70	+	4.7	+	1.4	+	9.99	C50
Butanol, t-	tert-Butanol, t-Buty alcohol	75-65-0	C <sub>4</sub> H <sub>10</sub> O	6.9	+	2.9	+			9.90	100
Butene, 1-	1-Butylene	106-98-9	C <sub>4</sub> H <sub>8</sub>			0.9				9.58	ne
Butoxyethanol, 2-	Butyl Cellosolve, Ethylene glycol monobutyl ether	111-76-2	C <sub>6</sub> H <sub>14</sub> O <sub>2</sub>	1.8	+	1.2	+	0.6	+	<10	25
Butyl acetate, n-		123-86-4	C <sub>6</sub> H <sub>12</sub> O <sub>2</sub>			2.6	+			10	150
Butyl acrylate, n-	Butyl 2-propenoate, Acrylic acid butyl ester	141-32-2	C <sub>7</sub> H <sub>12</sub> O <sub>2</sub>			1.6	+	0.6	+		10
Butylamine, n-		109-73-9	C <sub>4</sub> H <sub>11</sub> N	1.1	+	1.1	+	0.7	+	8.71	C5
Butyl cellosolve	see 2-Butoxyethanol	111-76-2									
Butyl hydroperoxide, t-		75-91-2	C <sub>4</sub> H <sub>10</sub> O <sub>2</sub>	2.0	+	1.6	+			<10	1
Butyl mercaptan	1-Butanethiol	109-79-5	C <sub>4</sub> H <sub>10</sub> S	0.55	+	0.52	+			9.14	0.5
Carbon disulfide		75-15-0	CS <sub>2</sub>	4	+	1.2	+	0.44		10.07	10
Carbon tetrachloride	Tetrachloromethane	56-23-5	CCl <sub>4</sub>	NR	+	NR	+	1.7	+	11.47	5
Carbonyl sulfide	Carbon oxysulfide	463-58-1	COS							11.18	
Cellosolve	see 2-Ethoxyethanol										
CFC-14	see Tetrafluoromethane										
CFC-113	see 1,1,2-Trichloro-1,2,2-trifluoroethane										
Chlorine		7782-50-5	Cl <sub>2</sub>					1.0	+	11.48	0.5
Chlorine dioxide		10049-04-4	ClO <sub>2</sub>	NR	+	NR	+	NR		10.57	0.1
Chloro-1,3-butadiene, 2-	Chloroprene	126-99-8	C <sub>4</sub> H <sub>5</sub> Cl			3					10

Compound Name	Synonym/Abbreviation	CAS No.	Formula	9.8	C	10.6	C	11.7	C	IP (eV)	TWA
Chlorobenzene	Monochlorobenzene	108-90-7	C <sub>6</sub> H <sub>5</sub> Cl	0.44	+	0.40	+	0.39	+	9.06	10
Chloro-1,1-difluoroethane, 1,1-	(R-142B)	75-68-3	C <sub>2</sub> H <sub>3</sub> ClF <sub>2</sub>			NR		NR		12.0	ne
Chlorodifluoromethane	HCFC-22, R-22	75-45-6	CHClF <sub>2</sub>	NR		NR		NR		12.2	1000
Chloroethane	Ethyl chloride	75-00-3	C <sub>2</sub> H <sub>5</sub> Cl	NR	+	NR	+	1.1	+	10.97	100
Chloroethanol	Ethylene chlorhydrin	107-07-3	C <sub>2</sub> H <sub>5</sub> ClO							10.52	C1
Chloroethyl ether, 2-	bis(2-chloroethyl) ether	111-44-4	C <sub>4</sub> H <sub>8</sub> Cl <sub>2</sub> O	8.6	+	3.0	+				5
Chloroethyl methyl ether, 2-	Methyl 2-chloroethyl ether	627-42-9	C <sub>3</sub> H <sub>7</sub> ClO			3					ne
Chloroform	Trichloromethane	67-66-3	CHCl <sub>3</sub>	NR	+	NR	+	3.5	+	11.37	10
Chloropicrin			CCl <sub>3</sub> NO <sub>2</sub>	NR	+	~400	+	7	+	?	0.1
Chlorotoluene, o-	o-Chloromethylbenzene	95-49-8	C <sub>7</sub> H <sub>7</sub> Cl			0.5		0.6		8.83	50
Chlorotoluene, p-	p-Chloromethylbenzene	106-43-4	C <sub>7</sub> H <sub>7</sub> Cl					0.6		8.69	ne
Crotonaldehyde	<i>trans</i> -2-Butenal	123-73-9	C <sub>4</sub> H <sub>6</sub> O	1.5	+	1.1	+	1.0	+	9.73	2
		4170-30-3									
Cumene	Isopropylbenzene	98-82-8	C <sub>9</sub> H <sub>12</sub>	0.58	+	0.54	+	0.4	+	8.73	50
Cyanogen bromide		506-68-3	CNBr	NR		NR		NR		11.84	ne
Cyanogen chloride		506-77-4	CNCl	NR		NR		NR		12.34	C0.3
Cyclohexane		110-82-7	C <sub>6</sub> H <sub>12</sub>	3.3	+	1.4	+	0.64	+	9.86	300
Cyclohexanol	Cyclohexyl alcohol	108-93-0	C <sub>6</sub> H <sub>12</sub> O					1.1		9.75	50
Cyclohexanone		108-94-1	C <sub>6</sub> H <sub>10</sub> O	1.0	+	0.9	+	0.7	+	9.14	25
Cyclohexene		110-83-8	C <sub>6</sub> H <sub>10</sub>			0.8	+			8.95	300
Cyclohexylamine		108-91-8	C <sub>6</sub> H <sub>13</sub> N			1.2				8.62	10
Cyclopentane		287-92-3	C <sub>5</sub> H <sub>10</sub>					0.6		10.51	600
Decane		124-18-5	C <sub>10</sub> H <sub>22</sub>	4.0	+	1.4	+	0.35	+	9.65	ne
Diacetone alcohol	4-Methyl-4-hydroxy-2-pentanone	123-42-2	C <sub>6</sub> H <sub>12</sub> O <sub>2</sub>			0.7					50
Dibromochloromethane	Chlorodibromomethane	124-48-1	CHBr <sub>2</sub> Cl	NR	+	5.3	+	0.7	+	10.59	ne
Dibromoethane, 1,2-	EDB, Ethylene dibromide, Ethylene bromide	106-93-4	C <sub>2</sub> H <sub>4</sub> Br <sub>2</sub>	NR	+	1.7	+	0.6	+	10.37	ne
Dichlorobenzene, o-	1,2-Dichlorobenzene	95-50-1	C <sub>6</sub> H <sub>4</sub> Cl <sub>2</sub>	0.54	+	0.47	+	0.38	+	9.08	25
Dichlorodifluoromethane	CFC-12	75-71-8	CCl <sub>2</sub> F <sub>2</sub>			NR	+	NR	+	11.75	1000
Dichloroethane, 1,2-	EDC, 1,2-DCA, Ethylene dichloride	107-06-2	C <sub>2</sub> H <sub>4</sub> Cl <sub>2</sub>			NR	+	0.6	+	11.04	10
Dichloroethene, 1,1-	1,1-DCE, Vinylidene chloride	75-35-4	C <sub>2</sub> H <sub>2</sub> Cl <sub>2</sub>			0.9				9.79	5
Dichloroethene, c-1,2-	c-1,2-DCE, <i>cis</i> -Dichloroethylene	156-59-2	C <sub>2</sub> H <sub>2</sub> Cl <sub>2</sub>			0.8				9.66	200
Dichloroethene, t-1,2-	t-1,2-DCE, <i>trans</i> -Dichloroethylene	156-60-5	C <sub>2</sub> H <sub>2</sub> Cl <sub>2</sub>			0.45	+	0.34	+	9.65	200
Dichloro-1-fluoroethane, 1,1-	R-141B	1717-00-6	C <sub>2</sub> H <sub>3</sub> Cl <sub>2</sub> F	NR	+	NR	+	2.0	+		ne
Dichloromethane	see Methylene chloride										
Dichloropentafluoropropane	AK-225, mix of ~45% 3,3-dichloro-1,1,1,2,2-pentafluoropropane (HCFC-225ca) & ~55% 1,3-Dichloro-1,1,2,2,3-pentafluoropropane (HCFC-225cb)	442-56-0	C <sub>3</sub> HCl <sub>2</sub> F <sub>5</sub>	NR	+	NR	+	25	+		ne
		507-55-1									
Dichloropropane, 1,2-		78-87-5	C <sub>3</sub> H <sub>6</sub> Cl <sub>2</sub>					0.7		10.87	75
Dichloro-1-propene, 1,3-		542-75-6	C <sub>3</sub> H <sub>4</sub> Cl <sub>2</sub>	1.3	+	0.96	+			<10	1
Dichloro-1-propene, 2,3-		78-88-6	C <sub>3</sub> H <sub>4</sub> Cl <sub>2</sub>	1.9	+	1.3	+	0.7	+	<10	ne
Dichloro-1,1,1-trifluoroethane, 2,2-	R-123	306-83-2	C <sub>2</sub> HCl <sub>2</sub> F <sub>3</sub>	NR	+	NR	+	10.1	+	11.5	ne
Dichlorvos	Vapona; O,O-dimethyl O-dichlorovinyl phosphate	62-73-7	C <sub>4</sub> H <sub>7</sub> Cl <sub>2</sub> O <sub>4</sub> P			0.9	+			<9.4	0.1
Dicyclopentadiene	DCPD, Cyclopentadiene dimer	77-73-6	C <sub>10</sub> H <sub>12</sub>	0.57	+	0.48	+	0.43	+	8.8	5
Diesel Fuel #1		68334-30-5	m.w. 226			0.9	+				11
Diesel Fuel #2		68334-30-5	m.w. 216			0.7	+	0.4	+		11
Diethylamine		109-89-7	C <sub>4</sub> H <sub>11</sub> N			1	+			8.01	5
Diethylaminopropylamine, 3-		104-78-9	C <sub>7</sub> H <sub>18</sub> N <sub>2</sub>			1.3					ne

Compound Name	Synonym/Abbreviation	CAS No.	Formula	9.8	C	10.6	C	11.7	C	IP (eV)	TWA
Diethylmaleate		141-05-9	C <sub>8</sub> H <sub>12</sub> O <sub>4</sub>			4					ne
Diethyl sulfide	see Ethyl sulfide										
Diisopropylamine		108-18-9	C <sub>6</sub> H <sub>15</sub> N	0.84	+	0.74	+	0.5	+	7.73	5
Diketene	Ketene dimer	674-82-8	C <sub>4</sub> H <sub>4</sub> O <sub>2</sub>	2.6	+	2.0	+	1.4	+	9.6	0.5
Dimethylacetamide, N,N-	DMA	127-19-5	C <sub>4</sub> H <sub>9</sub> NO	0.87	+	0.8	+	0.8	+	8.81	10
Dimethylamine		124-40-3	C <sub>2</sub> H <sub>7</sub> N			1.5				8.23	5
Dimethyl carbonate	Carbonic acid dimethyl ester	616-38-6	C <sub>3</sub> H <sub>6</sub> O <sub>3</sub>	NR	+	~70	+	1.7	+	~10.5	ne
Dimethyl disulfide	DMDS	624-92-0	C <sub>2</sub> H <sub>6</sub> S <sub>2</sub>	0.2	+	0.20	+	0.21	+	7.4	ne
Dimethylethylamine	DMEA	598-56-1	C <sub>4</sub> H <sub>11</sub> N	1.1	+	1.0	+	0.9	+	7.74	~3
Dimethylformamide, N,N-	DMF	68-12-2	C <sub>3</sub> H <sub>7</sub> NO			0.8				9.13	10
Dimethylhydrazine, 1,1-	UDMH	57-14-7	C <sub>2</sub> H <sub>8</sub> N <sub>2</sub>			0.8	+	0.8	+	7.28	0.01
Dimethyl methylphosphonate	DMMP, methyl phosphonic acid dimethyl ester	756-79-6	C <sub>3</sub> H <sub>9</sub> O <sub>3</sub> P	NR	+	4.3	+	0.74	+	10.0	ne
Dimethyl sulfide		77-78-1	C <sub>2</sub> H <sub>6</sub> O <sub>4</sub> S	~23		~20	+	2.3	+		0.1
Dimethyl sulfide	see Methyl sulfide										
Dimethyl sulfoxide	DMSO, Methyl sulfoxide	67-68-5	C <sub>2</sub> H <sub>6</sub> OS			1.4	+			9.10	ne
Dioxane, 1,4-		123-91-1	C <sub>4</sub> H <sub>8</sub> O <sub>2</sub>			1.3				9.19	25
Dowtherm A	see Therminol										
DS-108F Wipe Solvent	Ethyl lactate/Isopar H/ Propoxypropanol ~7:2:1	97-64-3 64742-48-9 1569-01-3	m.w. 118	3.3	+	1.6	+	0.7	+		ne
Epichlorohydrin	ECH Chloromethyloxirane, 1-chloro2,3-epoxypropane	106-89-8	C <sub>2</sub> H <sub>5</sub> ClO	~200	+	8.5	+	1.4	+	10.2	0.5
Ethane		74-84-0	C <sub>2</sub> H <sub>6</sub>			NR	+	15	+	11.52	ne
Ethanol	Ethyl alcohol	64-17-5	C <sub>2</sub> H <sub>6</sub> O			12	+	8		10.47	1000
Ethanolamine	MEA, Monoethanolamine	141-43-5	C <sub>2</sub> H <sub>7</sub> NO	5.6	+	1.6	+			8.96	3
Ethene	Ethylene	74-85-1	C <sub>2</sub> H <sub>4</sub>			10	+	3		10.51	ne
Ethoxyethanol, 2-	Ethyl cellosolve, Ethylene glycol monoethyl ether	110-80-5	C <sub>4</sub> H <sub>10</sub> O <sub>2</sub>			1.3				9.6	5
Ethyl acetate		141-78-6	C <sub>4</sub> H <sub>8</sub> O <sub>2</sub>			4.6	+			10.01	400
Ethyl acrylate		140-88-5	C <sub>5</sub> H <sub>8</sub> O <sub>2</sub>			2.4	+	1.0	+	<10.3	5
Ethylamine		75-04-7	C <sub>2</sub> H <sub>7</sub> N			0.8				8.86	5
Ethylbenzene		100-41-4	C <sub>8</sub> H <sub>10</sub>	0.52	+	0.52	+	0.51	+	8.77	100
Ethylene glycol	1,2-Ethanediol	107-21-1	C <sub>2</sub> H <sub>6</sub> O <sub>2</sub>			16	+	6	+	10.16	C100
Ethylene oxide	Oxirane, Epoxyethane	75-21-8	C <sub>2</sub> H <sub>4</sub> O			13	+	3.5	+	10.57	1
Ethyl ether	Diethyl ether	60-29-7	C <sub>4</sub> H <sub>10</sub> O			1.1	+			9.51	400
Ethyl 3-ethoxypropionate	EEP	763-69-9	C <sub>7</sub> H <sub>14</sub> O <sub>3</sub>	1.2	+	0.75	+				ne
Ethyl formate		109-94-4	C <sub>3</sub> H <sub>6</sub> O <sub>2</sub>					1.9		10.61	100
Ethyl hexyl acrylate, 2-	Acrylic acid 2-ethylhexyl ester	103-11-7	C <sub>11</sub> H <sub>20</sub> O <sub>2</sub>			1.1	+	0.5	+		ne
Ethyl (S)-(-)-lactate see also DS-108F	Ethyl lactate, Ethyl (S)-(-)-hydroxypropionate	687-47-8 97-64-3	C <sub>5</sub> H <sub>10</sub> O <sub>3</sub>	13	+	3.2	+	1.6	+	~10	ne
Ethyl mercaptan	Ethanethiol	75-08-1	C <sub>2</sub> H <sub>6</sub> S	0.60	+	0.56	+			9.29	0.5
Ethyl sulfide	Diethyl sulfide	352-93-2	C <sub>4</sub> H <sub>10</sub> S			0.5	+			8.43	ne
Formaldehyde	Formalin	50-00-0	CH <sub>2</sub> O					0.6		10.87	C0.3
Formic acid		64-18-6	CH <sub>2</sub> O <sub>2</sub>	NR	+	NR	+	9	+	11.33	5
Furfural	2-Furaldehyde	98-01-1	C <sub>5</sub> H <sub>4</sub> O <sub>2</sub>			0.92	+	0.8	+	9.21	2
Furfuryl alcohol		98-00-0	C <sub>5</sub> H <sub>6</sub> O <sub>2</sub>			0.80	+			<9.5	10
Gasoline #1		8006-61-9	m.w. 72			0.9	+				300
Gasoline #2, 92 octane		8006-61-9	m.w. 93	1.3	+	1.0	+	0.5	+		300
Glutaraldehyde	1,5-Pentanedial, Glutaric dialdehyde	111-30-8	C <sub>5</sub> H <sub>8</sub> O <sub>2</sub>	1.1	+	0.8	+	0.6	+		C0.0 5
Halothane	2-Bromo-2-chloro-1,1,1-trifluoroethane	151-67-7	C <sub>2</sub> HBrClF <sub>3</sub>					0.6		11.0	50
HCFC-22	see Chlorodifluoromethane										
HCFC-123	see 2,2-Dichloro-1,1,1-trifluoroethane										
HCFC-141B	see 1,1-Dichloro-1-fluoroethane										
HCFC-142B	see 1-Chloro-1,1-difluoroethane										
HCFC-134A	see 1,1,1,2-Tetrafluoroethane										

Compound Name	Synonym/Abbreviation	CAS No.	Formula	9.8	C	10.6	C	11.7	C	IP (eV)	TWA
HCFC-225	see Dichloropentafluoropropane										
Heptane, n-		142-82-5	C <sub>7</sub> H <sub>16</sub>	45	+	2.8	+	0.60	+	9.92	400
Hexamethyldisilazane, 1,1,1,3,3,3-	HMDS	999-97-3	C <sub>6</sub> H <sub>19</sub> NSi <sub>2</sub>			0.2	+	0.2	+	~8.6	
Hexane, n-		110-54-3	C <sub>6</sub> H <sub>14</sub>	350	+	4.3	+	0.54	+	10.13	50
Hexanol, 1-	Hexyl alcohol	111-27-3	C <sub>6</sub> H <sub>14</sub> O	9	+	2.5	+	0.55	+	9.89	ne
Hexene, 1-		592-41-6	C <sub>6</sub> H <sub>12</sub>			0.8				9.44	30
Hydrazine		302-01-2	H <sub>4</sub> N <sub>2</sub>	>8	+	2.6	+	2.1	+	8.1	0.01
Hydrogen	Synthesis gas	1333-74-0	H <sub>2</sub>	NR	+	NR	+	NR	+	15.43	ne
Hydrogen cyanide	Hydrocyanic acid	74-90-8	HCN	NR	+	NR	+	NR	+	13.6	C4.7
Hydrogen peroxide		7722-84-1	H <sub>2</sub> O <sub>2</sub>	NR	+	NR	+	NR	+	10.54	1
Hydrogen sulfide		7783-06-4	H <sub>2</sub> S	NR	+	3.3	+	1.5	+	10.45	10
Iodine		7553-56-2	I <sub>2</sub>	0.1	+	0.1	+	0.1	+	9.40	C0.1
Iodomethane	Methyl iodide	74-88-4	CH <sub>3</sub> I	0.21	+	0.22	+	0.26	+	9.54	2
Isoamyl acetate	Isopentyl acetate	123-92-2	C <sub>7</sub> H <sub>14</sub> O <sub>2</sub>	10.1		2.1		1.0		<10	100
Isobutane	2-Methylpropane	75-28-5	C <sub>4</sub> H <sub>10</sub>			100	+	1.2	+	10.57	ne
Isobutanol	2-Methyl-1-propanol	78-83-1	C <sub>4</sub> H <sub>10</sub> O	19	+	3.8	+	1.5		10.02	50
Isobutene	Isobutylene, Methyl butene	115-11-7	C <sub>4</sub> H <sub>8</sub>	1.00	+	1.00	+	1.00	+	9.24	ne
Isobutyl acetate		110-19-0	C <sub>6</sub> H <sub>12</sub> O <sub>2</sub>			2.6					150
Isobutyl acrylate	Isobutyl 2-propenoate, Acrylic acid Isobutyl ester	106-63-8	C <sub>7</sub> H <sub>12</sub> O <sub>2</sub>			1.5	+	0.60	+		ne
Isoflurane	1-Chloro-2,2,2-trifluoroethyl difluoromethyl ether, forane	26675-46-7	C <sub>3</sub> H <sub>2</sub> ClF <sub>5</sub> O							~11.7	ne
Isooctane	2,2,4-Trimethylpentane	540-84-1	C <sub>8</sub> H <sub>18</sub>			1.2				9.86	ne
Isopar E Solvent	Isoparaffinic hydrocarbons	64741-66-8	m.w.121	1.7	+	0.8	+				ne
Isopar G Solvent	Photocopier diluent	64742-48-9	m.w. 148			0.8	+				ne
Isopar K Solvent	Isoparaffinic hydrocarbons	64742-48-9	m.w. 156	0.9	+	0.5	+	0.27	+		ne
Isopar L Solvent	Isoparaffinic hydrocarbons	64742-48-9	m.w. 163	0.9	+	0.5	+	0.28	+		ne
Isopar M Solvent	Isoparaffinic hydrocarbons	64742-47-8	m.w. 191			0.7	+	0.4	+		ne
Isopentane	2-Methylbutane	78-78-4	C <sub>5</sub> H <sub>12</sub>			8.2					ne
Isophorone		78-59-1	C <sub>9</sub> H <sub>14</sub> O					3		9.07	C5
Isoprene	2-Methyl-1,3-butadiene	78-79-5	C <sub>5</sub> H <sub>8</sub>	0.69	+	0.63	+	0.60	+	8.85	ne
Isopropanol	Isopropyl alcohol, 2-propanol	67-63-0	C <sub>3</sub> H <sub>8</sub> O	500	+	6.0	+	2.7		10.12	400
Isopropyl acetate		108-21-4	C <sub>5</sub> H <sub>10</sub> O <sub>2</sub>			2.6				9.99	250
Isopropyl ether	Diisopropyl ether	108-20-3	C <sub>6</sub> H <sub>14</sub> O			0.8				9.20	250
Jet fuel JP-4	Jet B, Turbo B, Wide cut type aviation fuel	8008-20-6 + 64741-42-0	m.w. 115			1.0	+	0.4	+		ne
Jet fuel JP-5	Jet 5, Kerosene type aviation fuel	8008-20-6 + 64747-77-1	m.w. 167			0.6	+	0.5	+		15
Jet fuel JP-8	Jet A-1, Kerosene type aviation fuel	8008-20-6 + 64741-77-1	m.w. 165			0.6	+	0.3	+		15
Limonene, D-	(R)-(+)-Limonene	5989-27-5	C <sub>10</sub> H <sub>16</sub>			0.33	+			~8.2	ne
Kerosene C10-C16 petro.distillate	– see Jet Fuels	8008-20-6									
MDI	- see 4,4'-Methylenebis(phenylisocyanate)										
Mesitylene	1,3,5-Trimethylbenzene	108-67-8	C <sub>9</sub> H <sub>12</sub>	0.36	+	0.35	+	0.3	+	8.41	ne
Methane	Natural gas	74-82-8	CH <sub>4</sub>	NR	+	NR	+	NR	+	12.51	ne
Methanol	Methyl alcohol, carbinol	67-56-1	CH <sub>4</sub> O	NR	+	NR	+	2.5	+	10.85	200
Methoxyethanol, 2-	Methyl cellosolve, Ethylene glycol monomethyl ether	109-86-4	C <sub>3</sub> H <sub>8</sub> O <sub>2</sub>	4.8	+	2.4	+	1.4	+	10.1	5
Methoxyethoxyethanol, 2-	2-(2-Methoxyethoxy)ethanol Diethylene glycol monomethyl ether	111-77-3	C <sub>7</sub> H <sub>16</sub> O	2.3	+	1.2	+	0.9	+	<10	ne
Methyl acetate		79-20-9	C <sub>3</sub> H <sub>6</sub> O <sub>2</sub>	NR	+	6.6	+	1.4	+	10.27	200
Methyl acrylate	Methyl 2-propenoate, acrylic acid methyl ester	96-33-3	C <sub>4</sub> H <sub>6</sub> O <sub>2</sub>			3.7	+	1.2	+	(9.9)	2
Methylamine	Aminomethane	74-89-5	CH <sub>5</sub> N			1.2				8.97	5
Methyl bromide	Bromomethane	74-83-9	CH <sub>3</sub> Br	110	+	1.7	+	1.3	+	10.54	1

Compound Name	Synonym/Abbreviation	CAS No.	Formula	9.8	C	10.6	C	11.7	C	IP (eV)	TWA
Methyl t-butyl ether	MTBE, <i>tert</i> -Butyl methyl ether	1634-04-4	C <sub>5</sub> H <sub>12</sub> O			0.9	+			9.24	40
Methyl cellosolve	see 2-Methoxyethanol										
Methyl chloride	Chloromethane	74-87-3	CH <sub>3</sub> Cl	NR	+	NR	+	0.74	+	11.22	50
Methylcyclohexane		107-87-2	C <sub>7</sub> H <sub>14</sub>	1.6	+	0.97	+	0.53	+	9.64	400
Methylene bis(phenyl-isocyanate), 4,4'-	MDI, Mondur M		C <sub>15</sub> H <sub>10</sub> N <sub>2</sub> O <sub>2</sub>	Very slow ppb level response							0.005
Methylene chloride	Dichloromethane	75-09-2	CH <sub>2</sub> Cl <sub>2</sub>	NR	+	NR	+	0.89	+	11.32	25
Methyl ether	Dimethyl ether	115-10-6	C <sub>2</sub> H <sub>6</sub> O	4.8	+	3.1	+	2.5	+	10.03	ne
Methyl ethyl ketone	MEK, 2-Butanone	78-93-3		0.86	+	0.9	+	1.1	+	9.51	200
Methylhydrazine	Monomethylhydrazine, Hydrazomethane	60-34-4	C <sub>2</sub> H <sub>6</sub> N <sub>2</sub>	1.4	+	1.2	+	1.3	+	7.7	0.01
Methyl isobutyl ketone	MIBK, 4-Methyl-2-pentanone	108-10-1	C <sub>6</sub> H <sub>12</sub> O	0.9	+	0.8	+	0.6	+	9.30	50
Methyl isocyanate	CH <sub>3</sub> NCO	624-83-9	C <sub>2</sub> H <sub>3</sub> NO	NR	+	4.6	+	1.5		10.67	0.02
Methyl isothiocyanate	CH <sub>3</sub> NCS	551-61-6	C <sub>2</sub> H <sub>3</sub> NS	0.5	+	0.45	+	0.4	+	9.25	ne
Methyl mercaptan	Methanethiol	74-93-1	CH <sub>4</sub> S	0.65		0.54		0.66		9.44	0.5
Methyl methacrylate		80-62-6	C <sub>5</sub> H <sub>8</sub> O <sub>2</sub>	2.7	+	1.5	+	1.2	+	9.7	100
Methyl nonafluorobutyl ether	HFE-7100DL	163702-08-7, 163702-07-6	C <sub>5</sub> H <sub>3</sub> F <sub>9</sub> O			NR	+	~35	+		ne
Methyl-1,5-pentane-diamine, 2- (coats lamp)	Dytek-A amine, 2-Methyl pentamethylenediamine	15520-10-2	C <sub>6</sub> H <sub>16</sub> N <sub>2</sub>			~0.6	+			<9.0	ne
Methyl propyl ketone	MPK, 2-Pentanone	107-87-9	C <sub>5</sub> H <sub>12</sub> O			0.93	+	0.79	+	9.38	200
Methyl-2-pyrrolidinone, N-	NMP, N-Methylpyrrolidone, 1-Methyl-2-pyrrolidinone, 1-Methyl-2-pyrrolidone	872-50-4	C <sub>5</sub> H <sub>9</sub> NO	1.0	+	0.8	+	0.9	+	9.17	ne
Methyl salicylate	Methyl 2-hydroxybenzoate	119-36-8	C <sub>8</sub> H <sub>8</sub> O <sub>3</sub>	1.3		0.9		0.9		~9	ne
Methylstyrene, α-	2-Propenylbenzene	98-83-9	C <sub>9</sub> H <sub>10</sub>			0.5				8.18	50
Methyl sulfide	DMS, Dimethyl sulfide	75-18-3	C <sub>2</sub> H <sub>6</sub> S	0.49	+	0.44	+	0.46	+	8.69	ne
Mineral spirits	Stoddard Solvent, Varsol 1	8020-83-5, 8052-41-3, 68551-17-7	m.w. 144			0.7	+	0.39	+		100
Mineral Spirits - Viscor 120B	Calibration Fluid, b.p. 156-207°C	8052-41-3	m.w. 142	1.0	+	0.7	+	0.3	+		100
Mustard	HD, Bis(2-chloroethyl) sulfide	505-60-2, 39472-40-7, 68157-62-0	C <sub>4</sub> H <sub>8</sub> Cl <sub>2</sub> S			0.6					0.0005
Naphthalene	Mothballs	91-20-3	C <sub>10</sub> H <sub>8</sub>	0.45	+	0.42	+	0.40	+	8.13	10
Nitric oxide		10102-43-9	NO	~6		5.2	+	2.8	+	9.26	25
Nitrobenzene		98-95-3	C <sub>6</sub> H <sub>5</sub> NO <sub>2</sub>	2.6	+	1.9	+	1.6	+	9.81	1
Nitroethane		79-24-3	C <sub>2</sub> H <sub>5</sub> NO <sub>2</sub>					3		10.88	100
Nitrogen dioxide		10102-44-0	NO <sub>2</sub>	23	+	16	+	6	+	9.75	3
Nitromethane		75-52-5	CH <sub>3</sub> NO <sub>2</sub>					4		11.02	20
Nitropropane, 2-		79-46-9	C <sub>3</sub> H <sub>7</sub> NO <sub>2</sub>					2.6		10.71	10
Nonane		111-84-2	C <sub>9</sub> H <sub>20</sub>			1.4				9.72	200
Octane, n-		111-65-9	C <sub>8</sub> H <sub>18</sub>	13	+	1.8	+			9.82	300
Pentane		109-66-0	C <sub>5</sub> H <sub>12</sub>	80	+	8.4	+	0.7	+	10.35	600
Peracetic acid	Peroxyacetic acid, Acetyl hydroperoxide	79-21-0	C <sub>2</sub> H <sub>4</sub> O <sub>3</sub>	NR	+	NR	+	2.3	+		ne
Peracetic/Acetic acid mix	Peroxyacetic acid, Acetyl hydroperoxide	79-21-0	C <sub>2</sub> H <sub>4</sub> O <sub>3</sub>			50	+	2.5	+		ne
Perchloroethene	PCE, Perchloroethylene, Tetrachloroethylene	127-18-4	C <sub>2</sub> Cl <sub>4</sub>	0.69	+	0.57	+	0.31	+	9.32	25
PGME	Propylene glycol methyl ether, 1-Methoxy-2-propanol	107-98-2	C <sub>6</sub> H <sub>12</sub> O <sub>3</sub>	2.4	+	1.5	+	1.1	+		100
PGMEA	Propylene glycol methyl ether acetate, 1-Methoxy-2-acetoxypropane, 1-Methoxy-2-propanol acetate	108-65-6	C <sub>6</sub> H <sub>12</sub> O <sub>3</sub>	1.65	+	1.0	+	0.8	+		ne
Phenol	Hydroxybenzene	108-95-2	C <sub>6</sub> H <sub>6</sub> O	1.0	+	1.0	+	0.9	+	8.51	5
Phosgene	Dichlorocarbonyl	75-44-5	CCl <sub>2</sub> O			No data				11.2	0.1

Compound Name	Synonym/Abbreviation	CAS No.	Formula	9.8	C	10.6	C	11.7	C	IP (eV)	TWA
Phosphine		7803-51-2	PH <sub>3</sub>	28		3.9	+	1.1	+	9.87	0.3
Photocopier Toner	Isoparaffin mix					0.5	+	0.3	+		ne
Picoline, 3-	3-Methylpyridine	108-99-6	C <sub>6</sub> H <sub>7</sub> N			0.9				9.04	ne
Pinene, α-		2437-95-8	C <sub>10</sub> H <sub>16</sub>			0.31	+	0.47		8.07	ne
Pinene, β-		18172-67-3	C <sub>10</sub> H <sub>16</sub>	0.38	+	0.37	+	0.37	+	~8	100
Piperylene, isomer mix	1,3-Pentadiene	504-60-9	C <sub>5</sub> H <sub>8</sub>	0.76	+	0.69	+	0.64	+	8.6	100
Propane		74-98-6	C <sub>3</sub> H <sub>8</sub>			NR	+	1.8	+	10.95	2500
Propanol, n-	Propyl alcohol	71-23-8	C <sub>3</sub> H <sub>8</sub> O			5		1.7		10.22	200
Propene	Propylene	115-07-1	C <sub>3</sub> H <sub>6</sub>	1.5	+	1.4	+	1.6	+	9.73	ne
Propionaldehyde	Propanal	123-38-6	C <sub>3</sub> H <sub>6</sub> O			1.9				9.95	ne
Propyl acetate, n-		109-60-4	C <sub>5</sub> H <sub>10</sub> O <sub>2</sub>			3.5				10.04	200
Propylene carbonate		108-32-7	C <sub>4</sub> H <sub>6</sub> O <sub>3</sub>			62	+	1	+	10.5	ne
Propylene glycol	1,2-Propanediol	57-55-6	C <sub>3</sub> H <sub>8</sub> O <sub>2</sub>	18		5.5	+	1.6	+	<10.2	ne
Propylene oxide	Methyloxirane	75-56-9	C <sub>3</sub> H <sub>6</sub> O	~240		6.6	+	2.9	+	10.22	20
		16088-62-3									
		15448-47-2									
Propyleneimine	2-Methylaziridine	75-55-8	C <sub>3</sub> H <sub>7</sub> N	1.5	+	1.3	+	1.0	+	9.0	2
Propyl mercaptan, 2-	2-Propanethiol, Isopropyl mercaptan	75-33-2	C <sub>3</sub> H <sub>8</sub> S	0.64	+	0.66	+			9.15	ne
Pyridine		110-86-1	C <sub>5</sub> H <sub>5</sub> N	0.78	+	0.7	+	0.7	+	9.25	5
Pyrrolidine (coats lamp)	Azacyclohexane	123-75-1	C <sub>4</sub> H <sub>9</sub> N	2.1	+	1.3	+	1.6	+	~8.0	ne
RR7300 (PGME/PGMEA)	70:30 PGME:PGMEA (1-Methoxy-2-propanol:1-Methoxy-2-acetoxypropane)	107-98-2	C <sub>4</sub> H <sub>10</sub> O <sub>2</sub> / C <sub>6</sub> H <sub>12</sub> O <sub>3</sub>			1.4	+	1.0	+		ne
Sarin	GB, Isopropyl methylphosphonofluoridate	107-44-8 50642-23-4	C <sub>4</sub> H <sub>10</sub> FO <sub>2</sub> P			~3					
Stoddard Solvent - see Mineral Spirits		8020-83-5									
Styrene		100-42-5	C <sub>8</sub> H <sub>8</sub>	0.45	+	0.40	+	0.4	+	8.43	20
Sulfur dioxide		7446-09-5	SO <sub>2</sub>			NR	+	NR	+	12.32	2
Sulfur hexafluoride		2551-62-4	SF <sub>6</sub>	NR		NR		NR		15.3	1000
Sulfuryl fluoride	Vikane	2699-79-8	SO <sub>2</sub> F <sub>2</sub>	NR		NR		NR		13.0	5
Tabun	Ethyl N, N-dimethylphosphoramidocyanidate	77-81-6	C <sub>5</sub> H <sub>11</sub> N <sub>2</sub> O <sub>2</sub> P			0.8					15ppt
Tetrachloroethane, 1,1,1,2-		630-20-6	C <sub>2</sub> H <sub>2</sub> Cl <sub>4</sub>					1.3		~11.1	ne
Tetrachloroethane, 1,1,2,2-		79-34-5	C <sub>2</sub> H <sub>2</sub> Cl <sub>4</sub>	NR	+	NR	+	0.60	+	~11.1	1
Tetraethyllead	TEL	78-00-2	C <sub>8</sub> H <sub>20</sub> Pb	0.4		0.3		0.2		~11.1	0.008
Tetraethyl orthosilicate	Ethyl silicate, TEOS	78-10-4	C <sub>8</sub> H <sub>20</sub> O <sub>4</sub> Si			0.7	+	0.2	+	~9.8	10
Tetrafluoroethane, 1,1,1,2-	HFC-134A	811-97-2	C <sub>2</sub> H <sub>2</sub> F <sub>4</sub>			NR		NR			ne
Tetrafluoroethene	TFE, Tetrafluoroethylene, Perfluoroethylene	116-14-3	C <sub>2</sub> F <sub>4</sub>			~15				10.12	ne
Tetrafluoromethane	CFC-14, Carbon tetrafluoride	75-73-0	CF <sub>4</sub>			NR	+	NR	+	>15.3	ne
Tetrahydrofuran	THF	109-99-9	C <sub>4</sub> H <sub>8</sub> O	1.9	+	1.7	+	1.0	+	9.41	200
Tetramethyl orthosilicate	Methyl silicate, TMOS	681-84-5	C <sub>4</sub> H <sub>12</sub> O <sub>4</sub> Si	10	+	1.9	+			~10	1
Therminol VP-1	DowthrmA,3:1 Diphenyl oxide:	101-84-8	C <sub>12</sub> H <sub>10</sub> O	0.9	+	0.7	+				ne
	Biphenyl	92-52-4	C <sub>12</sub> H <sub>10</sub>								
Toluene	Methylbenzene	108-88-3	C <sub>7</sub> H <sub>8</sub>	0.54	+	0.50	+	0.51	+	8.82	50
Toluene-2,4-diisocyanate	TDI, 4-Methyl-1,3-phenylene-2,4-diisocyanate	584-84-9	C <sub>9</sub> H <sub>6</sub> N <sub>2</sub> O <sub>2</sub>	1.4	+	1.4	+	2.0	+		0.002
Trichlorobenzene, 1,2,4-	1,2,4-TCB	120-82-1	C <sub>6</sub> H <sub>3</sub> Cl <sub>3</sub>	0.7	+	0.46	+			9.04	C5
Trichloroethane, 1,1,1-	1,1,1-TCA, Methyl chloroform	71-55-6	C <sub>2</sub> H <sub>3</sub> Cl <sub>3</sub>			NR	+	1	+	11	350
Trichloroethane, 1,1,2-	1,1,2-TCA	79-00-5	C <sub>2</sub> H <sub>3</sub> Cl <sub>3</sub>	NR	+	NR	+	0.9	+	11.0	10
Trichloroethene	TCE, Trichloroethylene	79-01-6	C <sub>2</sub> HCl <sub>3</sub>	0.62	+	0.54	+	0.43	+	9.47	50
Trichlorotrifluoroethane, 1,1,2-	CFC-113	76-13-1	C <sub>2</sub> Cl <sub>3</sub> F <sub>3</sub>			NR		NR		11.99	1000
Triethylamine	TEA	121-44-8	C <sub>6</sub> H <sub>15</sub> N	0.95	+	0.9	+	0.65	+	7.3	1
Triethyl borate	TEB; Boric acid triethyl ester, Boron ethoxide	150-46-9	C <sub>6</sub> H <sub>15</sub> O <sub>3</sub> B			2.2	+	1.1	+	~10	ne
Triethyl phosphate	Ethyl phosphate	78-40-0	C <sub>6</sub> H <sub>15</sub> O <sub>4</sub> P	~50	+	3.1	+	0.60	+	9.79	ne
Trifluoroethane, 1,1,2-		430-66-0	C <sub>2</sub> H <sub>3</sub> F <sub>3</sub>					34		12.9	ne

Compound Name	Synonym/Abbreviation	CAS No.	Formula	9.8 C	10.6 C	11.7 C	IP (eV)	TWA
Trimethylamine		75-50-3	C <sub>3</sub> H <sub>9</sub> N		0.9		7.82	5
Trimethylbenzene, 1,3,5-	- see Mesitylene	108-67-8						25
Trimethyl borate	TMB; Boric acid trimethyl ester, Boron methoxide	121-43-7	C <sub>3</sub> H <sub>9</sub> O <sub>3</sub> B		5.1 +	1.2 +	10.1	ne
Trimethyl phosphate	Ethyl phosphate	512-56-1	C <sub>3</sub> H <sub>9</sub> O <sub>4</sub> P		8.0 +	1.3 +	9.99	ne
Turpentine	Pinenes (85%) + other diisoprenes	8006-64-2	C <sub>10</sub> H <sub>16</sub>	0.4 +	0.3 +		~8	100
Undecane		1120-21-4	C <sub>11</sub> H <sub>24</sub>		2		9.56	ne
Varsol – see Mineral Spirits								
Vinyl acetate		108-05-4	C <sub>4</sub> H <sub>6</sub> O <sub>2</sub>	1.5 +	1.2 +	1.0 +	9.19	10
Vinyl bromide	Bromoethylene	593-60-2	C <sub>2</sub> H <sub>3</sub> Br		0.4		9.80	5
Vinyl chloride	Chloroethylene, VCM	75-01-4	C <sub>2</sub> H <sub>3</sub> Cl		2.0 +	0.6 +	9.99	5
Vinylidene chloride - see 1,1-Dichloroethene								
Vinyl-2-pyrrolidinone, 1-	NVP, N-vinylpyrrolidone, 1-ethenyl-2-pyrrolidinone	88-12-0	C <sub>6</sub> H <sub>9</sub> NO	1.0 +	0.8 +	0.9 +		ne
Viscor 120B - see Mineral Spirits - Viscor 120B Calibration Fluid								
Xylene, m-	1,3-Dimethylbenzene	108-38-3	C <sub>8</sub> H <sub>10</sub>	0.50 +	0.43 +	0.40 +	8.56	100
Xylene, o-	1,2-Dimethylbenzene	95-47-6	C <sub>8</sub> H <sub>10</sub>	0.57 +	0.59 +	0.69	8.56	100
Xylene, p-	1,4-Dimethylbenzene	106-42-3	C <sub>8</sub> H <sub>10</sub>		0.45 +	0.62 +	8.44	100
None				1	1	1		
Undetectable				1E+6	1E+6	1E+6		

## Appendix I:

### Example of Automatic Calculation of Correction Factors, TLVs and Alarm Limits for Mixtures

(Calculations performed using Excel version of this database, available on request)

Compound	CF 9.8 eV	CF 10.6 eV	CF 11.7eV	Mol. Frac	Conc ppm	TLV ppm	STEL Ppm
Benzene	0.55	0.53	0.6	0.01	1	0.5	2.5
Toluene	0.54	0.5	0.51	0.06	10	50	150
Hexane, n-	300	4.3	0.54	0.06	10	50	150
Heptane, n-	45	2.8	0.6	0.28	50	400	500
Styrene	0.45	0.4	0.42	0.06	10	20	40
Acetone	1.2	1.1	1.4	0.28	50	750	1000
Isopropanol	500	6	2.7	0.28	50	400	500
None	1	1	1	0.00	0	1	
Mixture Value:	2.1	1.5	0.89	1.00	181	56	172
TLV Alarm Setpoint when Calibrated to Isobutylene:	26	37	62		ppm	ppm	ppm
	ppm	ppm	ppm				
STEL Alarm Setpoint, same Calibration	86	115	193				
	ppm	ppm	ppm				