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AnsellGUARDIAN® Chemical Report

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Disclaimer

In this report, you will find information related to the barrier performance of certain personal protective equipment (PPE) against the chemicals you selected. This information is intended to enable the Health and Safety professional at your organization make more informed decisions about the Ansell PPE that may offer the greatest protection in the intended circumstances and assist with carrying out a risk assessment for your organization.

We wish to highlight that permeation times do not equate to safe wear time. Safe wear time may vary depending on whether the PPE is donned correctly, the surrounding temperature, the chemicals' toxicity, and other factors. Permeation information offered here is limited to the main protective material. Permeation times may vary around seams, zips, visors or any other joins or components of the PPE. It is the responsibility of your organization's Health and Safety professional to undertake a risk assessment before choosing the appropriate PPE for the task at hand. If you want to discuss any aspect in detail, please contact us.

Estimations of the barrier properties of PPE are based on currently available data and extrapolations from laboratory test results and information regarding the chemicals' composition. Synergistic effects of mixing chemicals have not been accounted for. Estimations are subject to change if new testing is carried out or new information is available providing better grounds for extrapolations. For these reasons, any information in this report is provided for informational purposes only and Ansell fully disclaims any liability including warranties related to any statement contained herein.



Legend for Hand Protection

| Permeation Breakthrough Times | | | | | |
|-------------------------------|---------|-------------------|--|--|--|
| | <10 | Not Recommended | | | |
| | 10-30 | Splash Protection | | | |
| | 30-60 | Splash Protection | | | |
| | 60-120 | Medium Protection | | | |
| | 120-240 | Medium Protection | | | |
| | 240-480 | Good Protection | | | |
| | >480 | Good Protection | | | |

Permeation breakthrough time is the time (in minutes) for the chemical in question to be permeating through the material at a rate of 1.0 μg /cm²/min (as per EN ISO 374) or 0.1 μg /cm²/min (as per ASTM F739).

PS = Physical State: A = Aerosol, G = Gas, L = Liquid, P = Paste, S = Solid





Thickness (mm) : 0.35 mm / 13.8 mil

| CAS | Chemical Name | % | PS | EN ISO 374 | ASTM F739 |
|-----------|----------------------|-------|----|------------|-----------|
| 107-06-2 | 1,2-Dichloroethane | 100.0 | L | 18' C | |
| 95-49-8 | 2-Chlorotoluene | 100.0 | L | < 5' C | |
| 122-99-6 | 2-Phenoxyethanol | 100.0 | L | > 480' C | > 480' C |
| 106-95-6 | 3-Bromo-1-propene | 100.0 | L | 5' C | |
| 107-85-7 | 3-Methylbutylamin | 100.0 | L | < 10' C | |
| 64-19-7 | Acetic acid | 100.0 | L | > 480' C | |
| 67-64-1 | Acetone | 100.0 | L | 337' C | |
| 75-05-8 | Acetonitrile | 100.0 | L | > 480' C | |
| 106-92-3 | Allyl glycidyl ether | 100.0 | L | 335' C | |
| 7664-41-7 | Ammonia, gas | 100.0 | G | > 480' C | > 480' C |
| 62-53-3 | Aniline | 100.0 | L | > 480' C | > 480' C |
| 100-66-3 | Anisole | 100.0 | L | 21' C | 16' C |
| 71-43-2 | Benzene | 100.0 | L | 9' C | |
| 7726-95-6 | Bromine | 100.0 | L | 3' C | |
| 75-15-0 | Carbon disulfide | 100.0 | L | < 5' C | |
| 7782-50-5 | Chlorine, gas | 100.0 | G | > 480' C | > 480' C |
| 123-42-2 | Diacetone Alcohol | 100.0 | L | > 480' C | |
| 111-92-2 | Dibutylamine | 100.0 | L | < 10' C | < 10' C |
| 1300-21-6 | Dichloroethane | 100.0 | L | 18' C | |
| 75-09-2 | Dichloromethane | 100.0 | L | < 5' C | |





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| CAS | Chemical Name | % | PS | EN ISO 374 | ASTM F739 |
|------------|----------------------------------|-------|----|------------|-----------|
| 109-89-7 | Diethylamine | 100.0 | L | < 5' C | |
| 111-96-6 | Diethylene glycol dimethyl ether | 100.0 | L | 187' c | |
| 28454-70-8 | Diisononylamin | 100.0 | L | 12' C | |
| 108-20-3 | Diisopropylether | 100.0 | L | 8' C | 2' C |
| 927-62-8 | Dimethylbutylamine | 100.0 | L | < 10' C | < 10' C |
| 111-43-3 | Dipropyl ether | 100.0 | L | < 1' C | < 1' C |
| 106-89-8 | Epichlorohydrin | 100.0 | L | > 480' C | |
| 141-78-6 | Ethyl acetate | 100.0 | L | 55' C | |
| 105-54-4 | Ethyl butyrate | 100.0 | L | 17' C | 19' C |
| 110-71-4 | Ethylene Glycol Dimethyl Ether | 100.0 | L | < 1' C | 2' C |
| 50-00-0 | Formaldehyde | 37.0 | L | > 480' C | |
| 110-00-9 | Furan | 100.0 | L | 8' C | < 1' C |
| 142-82-5 | Heptane | 100.0 | L | 5' C | |
| 392-56-3 | Hexafluorobenzene | 100.0 | L | 77' c | |
| 7664-39-3 | Hydrofluoric Acid | 40.0 | L | > 480' C | |
| 7722-84-1 | Hydrogen peroxide | 30.0 | L | > 480' C | |
| 78-81-9 | Isobutylamine | 100.0 | L | < 10' C | |
| 27775-00-4 | Isononylamin | 100.0 | L | 12' c | |
| 67-56-1 | Methanol | 100.0 | L | > 480' C | |
| 78-93-3 | Methyl ethyl ketone | 100.0 | L | 76' c | |





Thickness (mm) : 0.35 mm / 13.8 mil

| CAS | Chemical Name | % | PS | EN ISO 374 | ASTM F739 |
|-----------|-------------------------|-------|----|------------|-----------|
| 108-10-1 | Methyl Isobutyl Ketone | 100.0 | L | 69' C | |
| 127-19-5 | N,N-Dimethylacetamide | 100.0 | L | > 480' C | |
| 121-69-7 | N,N-Dimethylbenzenamine | 100.0 | L | 46' C | 41' C |
| 123-86-4 | n-Butyl acetate | 100.0 | L | 39' C | |
| 109-73-9 | n-Butylamine | 100.0 | L | < 10' C | < 10' C |
| 1126-78-9 | N-Butylaniline | 100.0 | L | 195' C | 135' C |
| 110-68-9 | N-Butylmethylamine | 100.0 | L | < 10' C | < 10' C |
| 872-50-4 | N-Methyl-2-pyrrolidone | 100.0 | L | > 480' C | > 480' C |
| 100-61-8 | N-Methylaniline | 100.0 | L | > 480' C | 443' C |
| 7697-37-2 | Nitric acid | 70.0 | L | > 480' C | |
| 7697-37-2 | Nitric acid | 65.0 | L | > 480' C | |
| 111-86-4 | Octylamine | 100.0 | L | 12' C | 10' C |
| 7601-90-3 | Perchloric acid | 70.0 | L | > 480' C | |
| 307-34-6 | Perfluorooctane | 100.0 | L | > 480' C | |
| 108-95-2 | Phenol | 90.0 | L | > 480' C | |
| 75-56-9 | Propylene Oxide | 100.0 | L | 8' C | |
| 110-86-1 | Pyridine | 100.0 | L | 89' C | |
| 1310-73-2 | Sodium Hydroxide | 40.0 | L | > 480' C | |
| 7664-93-9 | Sulfuric acid | 96.0 | L | 277' C | |
| 1634-04-4 | Tert-Butyl Methyl Ether | 100.0 | L | 11' c | |





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| CAS | Chemical Name | % | PS | EN ISO 374 | ASTM F739 |
|----------|--|-------|----|--------------------|-----------|
| 127-18-4 | Tetrachloroethylene | 100.0 | L | 5' C | |
| 109-99-9 | Tetrahydrofuran | 100.0 | L | 5' C | |
| 108-88-3 | Toluene | 100.0 | L | < 5' C | |
| 102-82-9 | Tributylamine | 100.0 | L | 41' C | 32' C |
| 79-01-6 | Trichloroethylene | 100.0 | L | < 5' C | |
| 149-73-5 | Trimethoxymethane | 100.0 | L | > 480' C | |
| | Hydrogen Fluoride (CAS# 7664-39-3, 17 C) | | L | 270' c | |
| | Phenol (CAS#108-95-2, 45 C, molten) | | L | > 480' C | > 480' C |
| | Phenol (CAS#108-95-2, 70 C, molten) | | L | 25' C | 15' C |
| | Phenol 30%(CAS#108-95-2, at 70 C) | | L | 37' C | 24' C |
| | Phenol 50%(CAS#108-95-2, at 70 C) | | L | 38' <mark>C</mark> | 22' C |

