<table>
<thead>
<tr>
<th>Compound</th>
<th>SMILES code</th>
<th>$p_0^*$ (298.15 K) (Pa)</th>
<th>$\Delta H_{\text{vap}}$ (kJ mol$^{-1}$)</th>
<th>$H$ (mol atm$^{-1}$)</th>
</tr>
</thead>
<tbody>
<tr>
<td>LNISOOH</td>
<td>O=CC(O)C(C)(OO)CON(=O)=O*</td>
<td>$2.2 \times 10^{-4}$</td>
<td>122.7</td>
<td>$2.1 \times 10^5$</td>
</tr>
<tr>
<td></td>
<td>CC(O)(CON(=O)=O)C(=O)C=O</td>
<td>$3.8 \times 10^{-4}$</td>
<td></td>
<td>120.0</td>
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<tr>
<td>LISPOOHOOOH</td>
<td>OC(C)(COO)C(C)=O*</td>
<td>$3.8 \times 10^{-7}$</td>
<td>155.3</td>
<td>$2.0 \times 10^{16}$</td>
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<tr>
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<td>CC(CO)(COO)O)OO</td>
<td>$1.9 \times 10^{-7}$</td>
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<td>158.9</td>
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<tr>
<td>LC578OOH</td>
<td>OCC(C)(CO)=O*</td>
<td>$2.0 \times 10^{-4}$</td>
<td>123.2</td>
<td>$3.0 \times 10^{11}$</td>
</tr>
<tr>
<td></td>
<td>O=CC(O)C(C)(CO)OO</td>
<td>$2.0 \times 10^{-4}$</td>
<td></td>
<td>123.2</td>
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<tr>
<td>C59OOH</td>
<td>OCC(=O)C(C)(CO)OO*</td>
<td>$1.0 \times 10^{-4}$</td>
<td>125.0</td>
<td>$3.0 \times 10^{11}$</td>
</tr>
</tbody>
</table>