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- tetrahydro-4H-pyrido[1,2-a]pyrimidin-4-one
- 6αβ-Hydroxy-17α-ethynyl-13α-methyl-7,8,9,11,12,13,14,15,17-decahydro-6H-cyclopenta[a]phenanthrene
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USP Pharmaceutical Analytical Impurities Catalog

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<td>(Forny lamidinoo) (N'-)Beta-(D-)Ribofuranosyl urea)</td>
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<td>(Budesonide S11)-ene 25 mg (8b-(2-hydroxyacetamido) -8b-(2-hydroxyacetamido) phenyl)</td>
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<td>Clopidogrel Open Ring Methyl Ester Hydrochloride (10 mg, methyl 2-(2-chlorophenyl)-2-(2-thiophen-2-yl)ethanimidate, hydrochloride)</td>
<td>Pharmaceutical Analytical Impurity</td>
<td>F10840</td>
<td>T10108-18-4</td>
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<td>1A00720</td>
<td>Active</td>
<td>Fluoro-3-methyl-7-oxo-10-(4-methylpiperazin-1-yl)-2,3-dihydro-7H-pyrido[1,2,3-de]1,4-benzoxazine</td>
<td>Pharmaceutical Analytical Impurity</td>
<td>F17280</td>
<td>23551-25-9</td>
<td>N/A</td>
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<td>2932996560</td>
<td>VIAL</td>
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<td>F162L0</td>
<td>Active</td>
<td>Desiccation Acid (25 mg) (1TBA + l,2,3,4-3-ethyl-5-17B eicos-8,11-diole-9-1,4-diene-17-carboxylic Acid)</td>
<td>Analytical Impurity</td>
<td>E162L0-0</td>
<td>01006000-0</td>
<td>5026-</td>
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<td>3-(9H-Carbazol-9-yl) 25 mg</td>
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<td>0.77/0.00</td>
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<td>mg</td>
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<td>Cold Shipment Required</td>
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**Product Name:** Fluoro-11-hydroxy-10,13-trimethyl-3,6,7,8,9,10,11,12,13,14,15,16-dodecahydro-17H-cyclopenta[a]phenanthren-17-ylidene-3-hydroxyacetaldehyde (25 mg)

**CAS Number:** 37926-77-5

**NDC Number:** 54052-00

**Net Weight:** 25 mg

**Unit of Measure:** mg

**Commodity Codes (HS Codes):** 2011899400

**Special Restriction:** Cold Shipment Required

**Package Type:** VIAL

**USMCA Eligible:** No

**KORUS Eligible:** No

**Base Control Drug:** No

**Base Control Drug %:** No
| Catalog# | Status | Product Type | Current Lot | Previous Lot | CAS# | NDC# | Unit Price | Co. of Origin | Material Origin | UN # | Net Weight | Unit of Measure | Commodity Codes (HS Codes)* | Special Restriction | Pkg. Type | USMCA Eligible | KORUS Eligible | Base Control Drug | Base Control Drug % |
|----------|--------|--------------|-------------|--------------|------|------|------------|--------------|-----------------|------|-------------|-----------------|-------------------|-------------------|-----------|----------------|----------------|-------------------|----------------|-----------------|
| TA50020 | Active | Drug        |             |              |      |      |            |              |                 |      |             |                 |                   |                   |           |               |                |                   |                |                 |
|          |        |              |              |              |      |      |            |              |                 |      |             |                 |                   |                   |           |               |                |                   |                |                 |

**Entorexine (Ethyl 11Beta,17alpha-dihydroxy-5-fluro-16alpha-methyl andros-14-diene-3-one-17-ylcarboxylate)**

<p>| Catalog# | Status | Product Type | Current Lot | Previous Lot | CAS# | NDC# | Unit Price | Co. of Origin | Material Origin | UN # | Net Weight | Unit of Measure | Commodity Codes (HS Codes)* | Special Restriction | Pkg. Type | USMCA Eligible | KORUS Eligible | Base Control Drug | Base Control Drug % |
|----------|--------|--------------|-------------|--------------|------|------|------------|--------------|-----------------|------|-------------|-----------------|-------------------|-------------------|-----------|----------------|----------------|-------------------|----------------|-----------------|
| F169B0  |        | Active Drug  |             |              |      |      |            |              |                 |      |             |                 |                   |                   |           |               |                |                   |                |                 |
|          |        |              |              |              |      |      |            |              |                 |      |             |                 |                   |                   |           |               |                |                   |                |                 |</p>
<table>
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<th>Catalog #</th>
<th>Status</th>
<th>Product Name</th>
<th>Product Type</th>
<th>Current Lot</th>
<th>Previous Lot</th>
<th>CAS#</th>
<th>NDC#</th>
<th>Unit Price</th>
<th>Net Weight</th>
<th>Unit Of Measure</th>
<th>Commodity Codes (HS Codes)*</th>
<th>Special Restriction</th>
<th>Pkg. Type</th>
<th>USMCA Eligible</th>
<th>KORUS Eligible</th>
<th>Base Control Drug</th>
<th>Base Control Drug %</th>
</tr>
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<tbody>
<tr>
<td>1A000250</td>
<td>Active</td>
<td>(R)-9-fluoro-17-(formyl oxy)-11-hydroxy-10,13,16-trimethyl-3-oxo-6,7,8,9,12,13,14,15,16,17-dodecahydro-3H-cyclopenta[a]phenanthrene-17-carboxylic acid)</td>
<td>1A000250</td>
<td>844694-85-5</td>
<td>N/A</td>
<td>2019-00</td>
<td>Chem Synthesis</td>
<td>25</td>
<td>mg</td>
<td>23223396290</td>
<td>VAL</td>
<td>No</td>
<td>No</td>
<td></td>
<td></td>
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</tr>
</tbody>
</table>

USP Pharmaceutical Analytical Impurities Catalog

1A000250 Active (R)-9-fluoro-17-(formyl oxy)-11-hydroxy-10,13,16-trimethyl-3-oxo-6,7,8,9,12,13,14,15,16,17-dodecahydro-3H-cyclopenta[a]phenanthrene-17-carboxylic acid)
<table>
<thead>
<tr>
<th>Catalog #</th>
<th>Status</th>
<th>Product Name</th>
<th>CAS #</th>
<th>NDC #</th>
<th>Unit Price</th>
<th>Unit of Measure</th>
<th>Commodity Codes (HS Codes)*</th>
<th>Special Restriction</th>
<th>Plag Type</th>
<th>USMCA Eligible</th>
<th>KORUS Eligible</th>
<th>Base Control Drug</th>
<th>Base Control Drug %</th>
</tr>
</thead>
<tbody>
<tr>
<td>1A00360</td>
<td>Active</td>
<td>Bromide of (1,1-dibenzyl-4-(5,6-dimethoxy-1-oxo-2,3-dihydro-1H-inden-2-yl)methyl)piperidin-1-ium, bromide (25 mg)</td>
<td>F167Y0</td>
<td>86917-74-0</td>
<td>N/A</td>
<td>0.55</td>
<td>10</td>
<td>2933998350</td>
<td>VIAL</td>
<td>No</td>
<td>No</td>
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</tr>
<tr>
<td>1A00731</td>
<td>Active</td>
<td>Fexofenadine Olefin</td>
<td>F17290</td>
<td>1187954-57-9</td>
<td>N/A</td>
<td>0.56</td>
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<td>2933995070</td>
<td>VIAL</td>
<td>No</td>
<td>No</td>
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**Notes:**
- **Product Type:** Pharmaceutical Analytical Impurity
- **Material Origin:** CN Chemical Synthesis
- **Unit Of Measure:** VIAL
- **Special Restriction:** No
- **Base Control:** Drug
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<th>Catalog #</th>
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<th>Product Name</th>
<th>Product Type</th>
<th>Current Lot</th>
<th>Previous Lot</th>
<th>CAS #</th>
<th>NDC #</th>
<th>Unit Price</th>
<th>Unit of Measure</th>
<th>Commodity Codes (HS Codes)*</th>
<th>Special Restriction</th>
<th>Pkg Type</th>
<th>USMCA Eligible</th>
<th>KORUS Eligible</th>
<th>Base Control Drug</th>
<th>Base Control Drug %</th>
</tr>
</thead>
<tbody>
<tr>
<td>1AD0120</td>
<td>Active</td>
<td>Fluoro androstadiene Carboxylic Acid (25 mg)</td>
<td>Pharmaceutical Analytical Impurity</td>
<td>37927-01-8</td>
<td>34988-00</td>
<td>UNI</td>
<td>Chemical Synthesis</td>
<td>25</td>
<td>Tab</td>
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</table>

**Detail:**
- **Product Name:** Fluoro androstadiene Carboxylic Acid (25 mg)
- **CAS Number:** 37927-01-8
- **NDC Number:** 34988-00
- **Unit of Measure:** Tab
- **Commodity Codes (HS Codes):** 2918309900

**Formula:**
(2S, 4S, 4aS, 5aR, 6S, 12bR, 12cR, 12eR, 12fR, 12gR, 12hR)-4a-fluoro-5,12b,12c,12e,12f,12g,12h,12h-tetrahydroxychrysene (25 mg)
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<th>Catalog#</th>
<th>Status</th>
<th>Product Name</th>
<th>CAS#</th>
<th>NDC#</th>
<th>Co. of Origin</th>
<th>Material Origin</th>
<th>UN #</th>
<th>Net Weight</th>
<th>Unit Of Measure</th>
<th>Commodity Codes (HS Codes)*</th>
<th>Special Restriction</th>
<th>Pkg. Type</th>
<th>USMCA Eligible</th>
<th>KORUS Eligible</th>
<th>Base Control Drug</th>
<th>Base Control Drug %</th>
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</thead>
<tbody>
<tr>
<td>1A00310 A Active Hydroxy product of Acetophenone (25 mg)</td>
<td>16-trimethyl-3-oxo-6,7,8,9,10,11,12,13,14,15,16,17-dodecahydro-3H-cyclopenta[a]phenanthrene-17-carboxylic acid)</td>
<td>116814</td>
<td>1374817-71-5</td>
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<td>30</td>
<td>Tg</td>
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<td>Unit of Measure</td>
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<td>Pkg. Type</td>
<td>USMCA Eligible</td>
<td>KORUS Eligible</td>
<td>Base Control Drug</td>
<td>Base Control Drug %</td>
</tr>
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<td>Ethylene Glycol, Tetraethylene Glycol Diester (25 mg)</td>
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<td>2-Phenyl-1-propanone</td>
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<td>AMPULE</td>
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<td>1A00180</td>
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<td>Active N-Ethyl Quetiapine</td>
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<td>101158 03-4 (Free base)</td>
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<tr>
<td>1A00230</td>
<td>Active</td>
<td>N-[2-hydroxy-Carvool (8-di- (1-DMC Carbazol-4-syloxy)-3-(2-<a href="1-methyl-ethyl">2-(2-methoxyphenyl)ethyl</a>amino)-2-propanol)]hydrochloride</td>
<td>Pharmaceutical Analytical Impurity</td>
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<td>N</td>
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<tr>
<td>1A00530</td>
<td>Active</td>
<td>N6-[1-Iminoethyl]-D-Lysine ((2R)-2-<a href="1-methyl-ethyl">(2S)-2-amino-phenethyl</a>amino)-2-propanol)</td>
<td>Pharmaceutical Analytical Impurity</td>
<td>1844123-53-9</td>
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<td>N</td>
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<tr>
<td>1A00390</td>
<td>Active</td>
<td>Butanamide</td>
<td>116.2 6.21,4.5,6(6) 4.6-5 6,6-5 6,6-5 6,5-6 4,5-6 3,4-6 3,4-6 2,4-6</td>
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<tr>
<td>1A06646</td>
<td>Active</td>
<td>Penicillin V Ester (25 mg)</td>
<td>Dimethyl Ester (25 mg)</td>
<td>Parenteral: Intravenous Injection</td>
<td>Parenteral: Intravenous Injection</td>
<td>1554081-5</td>
<td>994</td>
<td>4600</td>
<td>VIAL</td>
<td>No No</td>
<td>No No</td>
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<td>Penicillin V Ester (25 mg)</td>
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<td>Commodity Codes (HS Codes)*</td>
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<td>KORUS Eligible</td>
<td>Base Control Drug</td>
<td>Base Control Drug %</td>
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<tr>
<td>1A0828</td>
<td>Water</td>
<td>Active Pemetr Disodium Enantiomer ((25 mg) (2R)-2-[[4-[2-(2-amino-4-oxo-4,7-dihydro-1H-pyrrolo[2,3-d]pyrimidin-5-yl)ethyl]benzoyl]amino]pentanedioic acid, sodium</td>
<td>F171L0 937370-10-0</td>
<td>NA</td>
<td>USP/00</td>
<td>Chemical Synthesis</td>
<td>St</td>
<td>mg</td>
<td>2903596855</td>
<td>VAL</td>
<td>No</td>
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<td>Product Type</td>
<td>Current Lot</td>
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<td>Unit Price</td>
<td>Unit Of Measure</td>
<td>Commodity Codes (HS Codes)*</td>
<td>Special Restriction</td>
<td>Pkg. Type</td>
<td>USMCA Eligible</td>
<td>KORUS Eligible</td>
<td>Base Control Drug</td>
<td>Base Control Drug %</td>
</tr>
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</tr>
<tr>
<td>1A00610</td>
<td>Active</td>
<td>Active Pemetrax Glutam ide Trisodium Salt (25 mg) (((S)-4-((2-(2-amino-4-oxo-4,7-dihydro-1H-pyrrolo[2,3-d]pyrimidin-5-yl)ethyl)benzamido)-4-carboxybutanoyl)-L-glutamic acid, sodium salt)</td>
<td>Pharmaceutical Analytical Impurity</td>
<td>F11740</td>
<td>TFA</td>
<td>500.00</td>
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<td>1g</td>
<td>2030309500</td>
<td>VAL</td>
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<tr>
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<td>Active Pyrazinediethanol</td>
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<td>Commodity Codes (HS Codes)*</td>
<td>Special Restriction</td>
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<td>1A00380</td>
<td>Active</td>
<td>Pyrazinediethanol Propionic Acid (5 mg) (3-(6-bis(2-hydroxyethyl)pyrazin-2-yl)propionate), sodium salt</td>
<td>Pharmaceutical Analytical Impurity</td>
<td>F167X0</td>
<td>96681-85-5</td>
<td>N/A</td>
<td>37.2</td>
<td>00.0</td>
<td>N/A</td>
<td>Chemical Synthesis</td>
<td>5</td>
<td>Fig</td>
<td>2321390670</td>
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<td>No</td>
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<tr>
<td>1A00460</td>
<td>Active</td>
<td>Pyrrolic acid Derivatives (10 mg) (4-(2-hydroxyethyl)-1H-pyrrole-3-carboxylic acid), sodium salt</td>
<td>Pharmaceutical Analytical Impurity</td>
<td>F16860</td>
<td>404839-11-8</td>
<td>N/A</td>
<td>81.2</td>
<td>00.0</td>
<td>N/A</td>
<td>Chemical Synthesis</td>
<td>10</td>
<td>Fig</td>
<td>2321390552</td>
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<td>Base Control Drug %</td>
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<td>1A00190</td>
<td>Active</td>
<td>Quetiapine Dihydrochloride (COLD SHIPMENT REQUIRED)</td>
<td>Pharmaceutical Analytical Impurity</td>
<td>R165N0</td>
<td>F160H0 (30-NOV-2022)</td>
<td>329216-67-3</td>
<td>N/A</td>
<td>N/A</td>
<td>Chemical Synthesis</td>
<td>(25 mg)</td>
<td>Cold Shipment Required</td>
<td>VIAL</td>
<td>No</td>
<td>No</td>
<td>Quetiapine Base Control Drug</td>
<td>10%</td>
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<tr>
<td>1A00130</td>
<td>Active</td>
<td>Quetiapine Tetraethylammonium Fumarate Salt (COLD SHIPMENT REQUIRED)</td>
<td>Pharmaceutical Analytical Impurity</td>
<td>F158K0</td>
<td>2934990002</td>
<td>25 mg</td>
<td>2934990002</td>
<td>Cold Shipment Required</td>
<td>VIAL</td>
<td>No</td>
<td>No</td>
<td>No</td>
<td>No</td>
<td>No</td>
<td>Quetiapine Base Control Drug</td>
<td>10%</td>
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<td>1A00140</td>
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<td>Ezetimibe (25 mg)</td>
<td>Pharmaceutical Analytical Impurity</td>
<td>F159N0 163380-16-3</td>
<td>N/A</td>
<td>16-3</td>
<td>TEN</td>
<td>Parenteral Synthesis</td>
<td>54</td>
<td>mg</td>
<td>XX</td>
<td>VIAL</td>
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<td>1A00370</td>
<td>Active</td>
<td>(1R)-N-(1-</td>
<td>Levitiracetam Hydroxybutyramide Analog (10 mg) (R)-N-1-aminomethyl-3-hydroxybutyramide</td>
<td>11287-9</td>
<td>TCA</td>
<td>$450.00</td>
<td>F170F0</td>
<td>1847568-87-9</td>
<td>Chemical Synthesis</td>
<td>CN Chemical Synthesis</td>
<td>10 mg</td>
<td>VIAL</td>
<td>No</td>
<td>No</td>
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<tr>
<td>1A00600</td>
<td>Active</td>
<td>(3R)-3-(2-(4-</td>
<td>Risperidone Difluoroketone (10 mg) (3-&lt;p&gt;(2-(4-difluorobenzoyl)piperdin-1-yl)ethyl)-2-methyl-6,7,8,9&lt;&lt;/p&gt;Pharmaceutical Analytical Impurity</td>
<td>118101-67-7</td>
<td>TCA</td>
<td>$350.00</td>
<td>F167Q0</td>
<td>158697-67-7</td>
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<td>CN Chemical Synthesis</td>
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<tr>
<td>1A00670</td>
<td>Active</td>
<td>N-(3-cyano-3-[(10-(4-(3-oxomorpholino)phenyl)propyl)thieno-2-carboxamide</td>
<td>F171C0</td>
<td>N/A</td>
<td>500.00</td>
<td>10</td>
<td>mg</td>
<td>COTTON SYNTHESIS</td>
<td>2020900002</td>
<td>N/A</td>
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<tr>
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<td>Active</td>
<td>4H-pyrido[1,2-a]pyrimidin-4-one)</td>
<td>F170K0</td>
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<td>500.00</td>
<td>10</td>
<td>mg</td>
<td>COTTON SYNTHESIS</td>
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<td>1A006001</td>
<td>Active</td>
<td>(S)-4-[(4-(2-hydroxy-3-(methylamino)propyl)amino)morpholin-3-yl]morpholine Phthalamide (25 mg)</td>
<td>Pharmaceutical Analytical Impurity</td>
<td>F170L0</td>
<td>1365267-36-2</td>
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<td>000000</td>
<td>0.00</td>
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<td>40</td>
<td>2934990002</td>
<td>VIAL</td>
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<tr>
<td>1A00170</td>
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<td>Rosuvastatin Isoamyl Ester (25 mg) ((3S,5R,6E)-7-[4-(4-flurophenyl)-6-(1-methylethyl)-2-[N-methyl(methylsulphonyl)amino]-5-pyrimidinyl]-3,5-dihydroxy-6-heptanoic acid, methyl butyl (isoamyl ester))</td>
<td>Pharmaceutical Analytical Impurity</td>
<td>F158U0</td>
<td>119-045-06-3</td>
<td>N/A</td>
<td>$650.00</td>
<td>K</td>
<td>Chemical Synthesis</td>
<td>25</td>
<td>mg</td>
<td>2920908550</td>
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<td>No</td>
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<tr>
<td>1A00050</td>
<td>Active</td>
<td>Rosuvastatin Ketone (25 mg) ((R,E)-7-[(4-flurophenyl)-6-(1-methylethyl)-2-[N-methyl(methylsulphonyl)amino]-5-pyrimidinyl]-3,5-dihydroxy-6-heptanoic acid, methyl butyl (keto ester))</td>
<td>Pharmaceutical Analytical Impurity</td>
<td>R170M0</td>
<td>F15970 (31-OCT-2022)</td>
<td>N/A</td>
<td>$795.00</td>
<td>Chemical Synthesis</td>
<td>25</td>
<td>mg</td>
<td>2920908550</td>
<td>VIAL</td>
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<td>Pharmaceutical Analog (10 mg)</td>
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<td>3908100002</td>
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<td>Levitiracetam Hydroxybutyramide Analog (10 mg)</td>
<td>S4089</td>
<td>00</td>
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<td>3908100002</td>
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<td>1A00400</td>
<td>Active</td>
<td>(N,N',N',N'-tetramethylethylene diamine) (25 mg) (N1,N1,N2,N2-tetramethylethylene diamine)</td>
<td>Pharmaceutical Analytical Impurity F15800</td>
<td>110-18-9</td>
<td>N/A</td>
<td>54-1</td>
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<td>Chemical Synthesis</td>
<td>UN2372</td>
<td>25 mg</td>
<td>2801230055</td>
<td>AMPULE</td>
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<td>1A00750</td>
<td>Active</td>
<td>(3-bromo-N,N-dimethylaniline-4-oxo-4-[(p-tolyl)butanamide] (25 mg)</td>
<td>Pharmaceutical Analytical Impurity F172A0</td>
<td>836627-56-6</td>
<td>N/A</td>
<td>57-79-00</td>
<td>UN</td>
<td>Chemical Synthesis</td>
<td>UN</td>
<td>25 mg</td>
<td>2804246000</td>
<td>VIAL</td>
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<td>o-Celecoxib</td>
<td>Pharmaceutical Analytical Impurity F159P0</td>
<td>119589-99-0</td>
<td>N/A</td>
<td>404-05-0</td>
<td>UN</td>
<td>Chemical Synthesis</td>
<td>UN</td>
<td>25 mg</td>
<td>2809020050</td>
<td>VIAL</td>
<td>No</td>
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<td>Pkg. Type</td>
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*Disclaimer: The Commodity Codes (HS codes) provided in this catalog are for information purposes only and are subject to change without notice. The exporter and/or importer of record is responsible for