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## How to Use

- **Searching:** Type keyword in search field at top of page. Search by all or part of a monograph title. For searches using multiple criteria, you will find items that match each of the specified criteria unless quotation marks are used.
  - For example, a search on Aminosalicylic Acid Tablets will result in anything that contains “Aminosalicylic” OR “Acid” OR “Tablets”
  - A search for “Aminosalicylic Acid Tablets” will result in anything that specifically contains “Aminosalicylic Acid Tablets”
- **Sorting:** Click on any column header title to sort alphabetically or chronologically in ascending or descending order. Note: the page load column is sorted alphabetically so that a number is ordered by first digit vs. by the actual number; thus, numbers will not always be in order.
  - For example, page 2178 will come before page 74 on a page sort.
- **Downloading:** You can download the Errata table in Comma-separated Value (.csv). The download will include the Errata that you have filtered on.
- **Importing:** You will need to import the file into Excel or Open Office with UTF-8 encoding, as opposed to simply opening it. To import, open Excel or Open Office and select import from the File drop-down. Depending on the version you are using, you should be presented with import formatting options to include UTF-8 as one of the first steps. Importing via UTF-8 should eliminate odd character conversions.

<a href="#">Monograph Title</a> <a href="#">Section</a>	<a href="#">Source</a> <a href="#">Publication</a>	<a href="#">Page Number</a>	<a href="#">Errata Post</a> <a href="#">Date</a> <a href="#">Sort</a> <a href="#">ascending</a>	<a href="#">Errata Official</a> <a href="#">Date</a>	<a href="#">Target Errata</a> <a href="#">Print Publication</a>	<a href="#">Target Online</a> <a href="#">Fix Publication</a>	Description
CEFTIOFUR H ASSAY/	USPNF Online	Online	30-Sep-2022	1-Oct-2022	NA	NA	In <i>Analysis</i> :

<a href="#">Monograph Title</a>	<a href="#">Section</a>	<a href="#">Source Publication</a>	<a href="#">Page Number</a>	<a href="#">Errata Post Date</a>	<a href="#">Sort ascending</a>	<a href="#">Errata Official Date</a>	<a href="#">Target Errata Print Publication</a>	<a href="#">Target Online Fix Publication</a>	Description
YDROCHLORI DE									Change Calculate the percentage of ceftiofur (C <sub>19</sub> H <sub>17</sub> N <sub>5</sub> O <sub>7</sub> S <sub>3</sub> ) in the portion of Ceftiofur Hydrochloride taken: Result = $(r_U/r_S) \times (C_S/C_U) \times P \times 100$ to: Calculate the quantity, in µg/mg, of ceftiofur (C <sub>19</sub> H <sub>17</sub> N <sub>5</sub> O <sub>7</sub> S <sub>3</sub> ) in the portion of Ceftiofur Hydrochloride taken: Result = $(r_U/r_S) \times (C_S/C_U) \times P$
POTASSIUM GLUCONATE	IM PUR ITIES/ <i>Reducing Substances</i>	USPNF Online	Online	26-Aug-2022		1-Sep-2022	NA	NA	In <i>Titrimetric system</i> : Change <b>Titrant:</b> Iodine <b>Back-titrant:</b> Sodium thiosulfate

<a href="#">Monograph Title</a>	<a href="#">Section</a>	<a href="#">Source Publication</a>	<a href="#">Page Number</a>	<a href="#">Errata Post Date</a>	<a href="#">Sort ascending</a>	<a href="#">Errata Official Date</a>	<a href="#">Target Errata Print Publication</a>	<a href="#">Target Online Fix Publication</a>	Description
0.1 N Potassium Permanganate VS	STANDARDIZATION	USPNF Online	Online	26-Aug-2022		1-Sep-2022	NA	NA	to: <b>Titrant:</b> 0.1 N Iodine VS <b>Back-titrant:</b> 0.1 N Sodium Thiosulfate VS In <i>Standardization with potentiometric endpoint.</i> Change N = g Na <sub>2</sub> C <sub>2</sub> O <sub>4</sub> /mL KMnO <sub>4</sub> solution × 0.06700 to: N = g Na <sub>2</sub> C <sub>2</sub> O <sub>4</sub> /mL KMnO <sub>4</sub> solution × 0.06700
DACARBAZINE IM FOR INJECTION	PURITIES/ <i>Organic Impurities</i>	USPNF Online	Online	26-Aug-2022		1-Sep-2022	NA	NA	In <i>System suitability/Suitability require measurements/Signal-to-noise ratio:</i> Change

<a href="#">Monograph Title</a> <a href="#">Section</a>	<a href="#">Source Publication</a>	<a href="#">Page Number</a>	<a href="#">Errata Post Date</a> <a href="#">Sort ascending</a>	<a href="#">Errata Official Date</a>	<a href="#">Target Errata Print Publication</a>	<a href="#">Target Online Fix Publication</a>	Description
PROMETHAZINE HYDROCHLORIDE TABLETS IM PURITIES/ <i>Organic Impurities</i>	<i>USPNF Online</i>	Online	26-Aug-2022	1-Sep-2022	NA	NA	NTL 10, to: NLT 10, Change <b>System suitability solution:</b> 5 µg/mL each of USP Promethazine Hydrochloride RS and USP Promethazine Related Compound B RS from the <i>Standard stock solution</i> and <i>System suitability stock solution</i> , respectively <b>Standard solution:</b> 5 µg/mL of USP Promethazine Hydrochloride RS from the <i>Standard stock solution</i>

<a href="#">Monograph Title</a> <a href="#">Section</a>	<a href="#">Source Publication</a>	<a href="#">Page Number</a>	<a href="#">Errata Post Date</a> <a href="#">Sort ascending</a>	<a href="#">Errata Official Date</a>	<a href="#">Target Errata Print Publication</a>	<a href="#">Target Online Fix Publication</a>	Description
							<p><b>Sensitivity solution:</b> 0.25 µg/mL of USP Promethazine Hydrochloride RS from the <i>Standard solution</i> to:</p> <p><b>System suitability solution:</b> 5 µg/mL each of USP Promethazine Hydrochloride RS and USP Promethazine Related Compound B RS from the <i>Standard stock solution</i> and <i>System suitability stock solution</i>, respectively, in <i>Diluent Standard solution</i>: 5</p>

<a href="#">Monograph Title</a>	<a href="#">Section</a>	<a href="#">Source Publication</a>	<a href="#">Page Number</a>	<a href="#">Errata Post Date</a>	<a href="#">Sort ascending</a>	<a href="#">Errata Official Date</a>	<a href="#">Target Errata Print Publication</a>	<a href="#">Target Online Fix Publication</a>	Description
METHYLENE BLUE	ADDITIONAL REQUIREMENT S/USP Reference Standards <11>	USPNF Online	Online	26-Aug-2022		1-Sep-2022	NA	USPNF 2023 Issue 2	<p>µg/mL of USP Promethazine Hydrochloride RS from the <i>Standard stock solution</i> in <i>Diluent</i></p> <p><b>Sensitivity solution:</b> 0.25 µg/mL of USP Promethazine Hydrochloride RS from the <i>Standard solution</i> in <i>Diluent</i></p> <p>In USP Azure B RS: Change 3-(Dimethylamino)-7-(methylamino)-phenothiazine-5-ium chloride. to: 3-(Dimethylamino)-7-(methylamino)phenothiazine-5-ium chloride.</p> <p>In <i>System suitability</i></p>
METHSUXIMIDE	IMPUR	USPNF Online	Online	26-Aug-2022		1-Sep-2022	NA	NA	In <i>System suitability</i>

<a href="#">Monograph Title</a>	<a href="#">Section</a>	<a href="#">Source Publication</a>	<a href="#">Page Number</a>	<a href="#">Errata Post Date</a>	<a href="#">Sort ascending</a>	<a href="#">Errata Official Date</a>	<a href="#">Target Errata Print Publication</a>	<a href="#">Target Online Fix Publication</a>	Description
									<i>ity/Column efficiency: Change NTL 5800 theoretical plates to: NLT 5800 theoretical plates</i>
PROCHLORPE RAZINE MALEATE TABLETS	PERFORMANCE TESTS/ <i>Dissolution</i> <711>	USPNF Online	Online	26-Aug-2022		1-Sep-2022	NA	NA	<i>In Analysis: Change Results = <math>(A_U/A_S) \times C_S \times D \times (1/L) \times V \times 100</math> <math>A_U</math> = absorbance of the <i>Sample solution</i> <math>A_S</math> = absorbance of the <i>Standard solution</i> <math>C_S</math> = concentration of USP Prochlorpe razine Maleate RS in the <i>Standard solution</i></i>

<a href="#">Monograph Title</a> <a href="#">Section</a>	<a href="#">Source Publication</a>	<a href="#">Page Number</a>	<a href="#">Errata Post Date</a> <a href="#">Sort ascending</a>	<a href="#">Errata Official Date</a>	<a href="#">Target Errata Print Publication</a>	<a href="#">Target Online Fix Publication</a>	Description
							<p><math>D</math> = dilution factor for <i>Sample solution</i>, if needed</p> <p><math>L</math> = label claim (mg/Tablet)</p> <p><math>V</math> = volume of <i>Medium</i>, 500 mL</p> <p>to:</p> $\text{Result} = (A_U/A_S) \times C_S \times V \times D \times (M_{r1}/M_{r2}) \times (1/L) \times 100$ <p><math>A_U</math> = absorbance of the <i>Sample solution</i></p> <p><math>A_S</math> = absorbance of the <i>Standard solution</i></p> <p><math>C_S</math> = concentration of USP Prochlorperazine Maleate RS in the <i>Standard solution</i></p> <p><math>V</math> = volume of</p>



<a href="#">Monograph Title</a>	<a href="#">Section</a>	<a href="#">Source Publication</a>	<a href="#">Page Number</a>	<a href="#">Errata Post Date</a>	<a href="#">Sort ascending</a>	<a href="#">Errata Official Date</a>	<a href="#">Target Errata Print Publication</a>	<a href="#">Target Online Fix Publication</a>	Description
RESIDUAL SOLVENTS	USP REFERENCE STANDARDS <11>	USPNF Online	Online	26-Aug-2022		1-Sep-2022	NA	NA	<p>Medium, 500 mL</p> <p><math>D</math> = dilution factor for the <i>Sample solution</i>, if needed</p> <p><math>M_{r1}</math> = molecular weight of prochlorperazine, 373.94</p> <p><math>M_{r2}</math> = molecular weight of prochlorperazine maleate, 606.09</p> <p><math>L</math> = label claim (mg/Tablet)</p> <p>Change USP Residual Solvent Class 2—Mixture C RS to: USP Residual Solvents Class 2—Mixture C RS</p>
IRINOTECAN HYDROCHLORIDE INJECTION	DEFINITION	USPNF Online	Online	26-Aug-2022		1-Sep-2022	NA	NA	<p>Change NTL 90.0% and NMT 110.0% to: NLT 90.0% and NMT 110.0%</p>

<a href="#">Monograph Title</a>	<a href="#">Section</a>	<a href="#">Source Publication</a>	<a href="#">Page Number</a>	<a href="#">Errata Post Date</a>	<a href="#">Sort ascending</a>	<a href="#">Errata Official Date</a>	<a href="#">Target Errata Print Publication</a>	<a href="#">Target Online Fix Publication</a>	Description
0.1 M ZINC SULFATE VS	STANDARDIZATION	USPNF Online	Online	26-Aug-2022		1-Dec-2022	NA	NA	<p>In <i>Standardization with visual end point</i>. Change M = mL edetate disodium x edetate disodium/mL ZnSO<sub>4</sub> to: M = mL edetate disodium x M edetate disodium/mL ZnSO<sub>4</sub> AND</p> <p>In <i>Standardization with potentiometric end point</i>. Change M = mL edetate disodium x edetate disodium/mL ZnSO<sub>4</sub> to: M = mL edetate disodium x M edetate</p>

<a href="#">Monograph Title</a>	<a href="#">Section</a>	<a href="#">Source Publication</a>	<a href="#">Page Number</a>	<a href="#">Errata Post Date</a>	<a href="#">Errata Official Date</a>	<a href="#">Target Errata Print Publication</a>	<a href="#">Target Online Fix Publication</a>	Description
ECONAZOLE NITRATE	ADDITIONAL REQUIREMENT S/USP Reference Standards <11>	USPNF Online	Online	26-Aug-2022	1-Sep-2022	NA	USPNF 2023 Issue 2	disodium/mL ZnSO <sub>4</sub> In USP Econazole Related Compound C RS: Change 1-(4-Chlorobenzyl)-3-{2-[(4-chlorobenzyl)oxy]-2-(2,4-dichlorophenyl)-1H-imidazol-3-ium nitrate (salt). C <sub>25</sub> H <sub>21</sub> Cl <sub>14</sub> N <sub>3</sub> O <sub>4</sub> 569.26 to: 1-(4-Chlorobenzyl)-3-{2-[(4-chlorobenzyl)oxy]-2-(2,4-dichlorophenyl)-1H-imidazol-3-ium chloride. C <sub>25</sub> H <sub>21</sub> Cl <sub>5</sub> N <sub>2</sub> O 542.71
CHROMATOGRAPHY	ADJUSTMENT OF CHROMATOGRAPHIC	USPNF Online	Online	26-Aug-2022	1-Dec-2022	NA	NA	In Liquid Chromatography: Isocratic

<a href="#">Monograph Title</a> <a href="#">Section</a>	<a href="#">Source</a> <a href="#">Publication</a>	<a href="#">Page Number</a>	<a href="#">Errata Post</a> <a href="#">Date</a> <a href="#">Sort</a> <a href="#">ascending</a>	<a href="#">Errata Official</a> <a href="#">Date</a>	<a href="#">Target Errata</a> <a href="#">Print Publication</a>	<a href="#">Target Online</a> <a href="#">Fix Publication</a>	Description
CONDITIONS							<i>Elution/Injection volume:</i> Change $L_2$ = internal diameter of the column used (mm) $dc_1$ = particle size indicated in the monograph ( $\mu\text{m}$ ) $dc_2$ = particle size of the column used ( $\mu\text{m}$ ) to: $L_2$ = new column length (mm) $dc_1$ = column internal diameter indicated in the monograph (mm) $dc_2$ = new column internal diameter (mm) Change <b>System suitability</b>
PROMETHAZINE HYDROCHLORIDE IM PUR ITIES/ <i>Organic</i>	USPNF Online	Online	26-Aug-2022	1-Sep-2022	NA	NA	

<a href="#">Monograph Title</a> <a href="#">Section</a>	<a href="#">Source Publication</a>	<a href="#">Page Number</a>	<a href="#">Errata Post Date</a> <a href="#">Sort ascending</a>	<a href="#">Errata Official Date</a>	<a href="#">Target Errata Print Publication</a>	<a href="#">Target Online Fix Publication</a>	Description
<i>Impurities</i>							<p><b>solution:</b> 5 µg/mL each of USP Promethazine Hydrochloride RS and USP Promethazine Related Compound B RS from the <i>Standard stock solution</i> and <i>System suitability stock solution</i>, respectively</p> <p><b>Standard solution:</b> 5 µg/mL of USP Promethazine Hydrochloride RS from the <i>Standard stock solution</i></p> <p><b>Sensitivity solution:</b> 0.25 µg/mL of USP Promethazine Hydrochloride RS from the</p>

<a href="#">Monograph Title</a> <a href="#">Section</a>	<a href="#">Source Publication</a>	<a href="#">Page Number</a>	<a href="#">Errata Post Date</a> <a href="#">Sort ascending</a>	<a href="#">Errata Official Date</a>	<a href="#">Target Errata Print Publication</a>	<a href="#">Target Online Fix Publication</a>	Description
							<p><i>Standard solution</i> to: <b>System suitability solution:</b> 5 µg/mL each of USP Promethazine Hydrochloride RS and USP Promethazine Related Compound B RS from the <i>Standard stock solution</i> and <i>System suitability stock solution</i>, respectively, in <i>Diluent</i></p> <p><b>Standard solution:</b> 5 µg/mL of USP Promethazine Hydrochloride RS from the <i>Standard stock solution</i> in</p>

<a href="#">Monograph Title</a>	<a href="#">Section</a>	<a href="#">Source Publication</a>	<a href="#">Page Number</a>	<a href="#">Errata Post Date</a>	<a href="#">Sort ascending</a>	<a href="#">Errata Official Date</a>	<a href="#">Target Errata Print Publication</a>	<a href="#">Target Online Fix Publication</a>	Description
METHYLENE BLUE	IMPURITIES/ <i>Organic Impurities</i>	USPNF Online	Online	26-Aug-2022		1-Sep-2022	NA	USPNF 2023 Issue 2	<p><i>Diluent Sensitivity solution:</i> 0.25 µg/mL of USP Promethazine Hydrochloride RS from the <i>Standard solution</i> in <i>Diluent</i></p> <p>In footnote a in <i>Table 2:</i> Change 3-(Dimethylamino)-7-(methylo)-phenothiazine-5-ium chloride. to: 3-(Dimethylamino)-7-(methylo)phenothiazine-5-ium chloride.</p> <p>In <i>Analysis:</i> Change <math>W_U</math> = weight of the Oral Suspension taken (mg) to:</p>
MELOXICAM ORAL SUSPENSION	PERFORMANCE TESTS/ <i>Dissolution &lt;711&gt;</i>	USPNF Online	Online	29-Jul-2022		1-Aug-2022	NA	NA	<p>In <i>Analysis:</i> Change <math>W_U</math> = weight of the Oral Suspension taken (mg) to:</p>

<a href="#">Monograph Title</a> <a href="#">Section</a>	<a href="#">Source Publication</a>	<a href="#">Page Number</a>	<a href="#">Errata Post Date</a> <a href="#">Sort ascending</a>	<a href="#">Errata Official Date</a>	<a href="#">Target Errata Print Publication</a>	<a href="#">Target Online Fix Publication</a>	Description
LATANOPROS ASSAY/ T Procedure	USPNF Online	Online	29-Jul-2022	1-Dec-2022	NA	NA	<p><math>W_U</math> = weight of the Oral Suspension taken (g)</p> <p>Change <b>Standard solution:</b> Transfer 2.0 mg/mL of USP Latanoprost RS into a suitable volumetric flask, dissolve in dehydrated alcohol equivalent to 20% of the final volume, and dilute with chromatographic hexane to volume.</p> <p><b>Sample solution:</b> Transfer 2.0 mg/mL of Latanoprost into a suitable volumetric flask, dissolve in</p>



<a href="#">Monograph Title</a> <a href="#">Section</a>	<a href="#">Source Publication</a>	<a href="#">Page Number</a>	<a href="#">Errata Post Date</a> <a href="#">Sort ascending</a>	<a href="#">Errata Official Date</a>	<a href="#">Target Errata Print Publication</a>	<a href="#">Target Online Fix Publication</a>	Description
							<p>dehydrated alcohol equivalent to 20% of the final volume, and dilute with chromatographic hexane to volume.</p> <p>to:</p> <p><b>Standard solution:</b> 2.0 mg/mL of USP Latanoprost RS prepared as follows.</p> <p>Transfer USP Latanoprost RS into a suitable volumetric flask, dissolve in dehydrated alcohol equivalent to 20% of the final volume, and dilute with chromatographic hexane to volume.</p> <p><b>Sample</b></p>

<a href="#">Monograph Title</a>	<a href="#">Section</a>	<a href="#">Source Publication</a>	<a href="#">Page Number</a>	<a href="#">Errata Post Date</a>	<a href="#">Sort ascending</a>	<a href="#">Errata Official Date</a>	<a href="#">Target Errata Print Publication</a>	<a href="#">Target Online Fix Publication</a>	Description
FORSKOHLII	COMPOSITION <i>/Content of Forskolin</i>	USPNF Online	Online	29-Jul-2022		1-Aug-2022	NA	NA	<p><b>solution:</b> 2.0 mg/mL of Latanoprost prepared as follows. Transfer Latanoprost into a suitable volumetric flask, dissolve in dehydrated alcohol equivalent to 20% of the final volume, and dilute with chromatographic hexane to volume.</p> <p>In <i>Chromatographic system:</i> Change Column: 4.6-mm x 25-cm; 5-µm, 100 Å to: Column: 4.6-mm x 25-cm; 5-µm,</p>

<a href="#">Monograph Title</a>	<a href="#">Section</a>	<a href="#">Source Publication</a>	<a href="#">Page Number</a>	<a href="#">Errata Post Date</a>	<a href="#">Sort ascending</a>	<a href="#">Errata Official Date</a>	<a href="#">Target Errata Print Publication</a>	<a href="#">Target Online Fix Publication</a>	Description
SUTURES--NE PROCEDURE EDLE ATTACHMENT		USPNF Online	Online	29-Jul-2022		1-Aug-2022	NA	NA	100 Å; packing L1 In <i>Removable Needle Attachment</i> . Change For USP sizes 5-0 through 2-0, to: For USP sizes 5-0 through 2, Change
BACLOFEN INJECTION	SPECIFIC TESTS	USPNF Online	Online	29-Jul-2022		1-Aug-2022	NA	NA	<ul style="list-style-type: none"> <li>• <b>Osmolality and Osmolarity</b> ?785?, <i>Osmolality</i>: 270–320 mOsm/kg to:</li> <li>• <b>Osmolality and Osmolarity</b> ?785? <b>Osmolality</b>: 270–320 mOsm/kg</li> </ul>
POWDERED FORSKOHLII	COMPOSITION <i>/Content of Forskolin</i>	USPNF Online	Online	29-Jul-2022		1-Aug-2022	NA	NA	In <i>Chromatographic system</i> :

<a href="#">Monograph Title</a>	<a href="#">Section</a>	<a href="#">Source Publication</a>	<a href="#">Page Number</a>	<a href="#">Errata Post Date</a>	<a href="#">Sort ascending</a>	<a href="#">Errata Official Date</a>	<a href="#">Target Errata Print Publication</a>	<a href="#">Target Online Fix Publication</a>	Description
POWDERED FORSKOHLII EXTRACT	COMPOSITION /Content of Forskolin	USPNF Online	Online	29-Jul-2022		1-Aug-2022	NA	NA	Change Column: 4.6-mm x 25-cm; 5-µm, 100 Å to: Column: 4.6-mm x 25-cm; 5-µm, 100 Å; packing L1 In Chromatographic system: Change Column: 4.6-mm x 25-cm; 5-µm, 100 Å to: Column: 4.6-mm x 25-cm; 5-µm, 100 Å; packing L1
AMMONIUM G LYCYRRHIZATE	CHEMICAL INFORMATION	USPNF Online	Online	24-Jun-2022		1-Jul-2022	NA	NA	Change 840.08 to: 839.97
MELENGESTROL ACETATE	CHEMICAL INFORMATION	USPNF Online	Online	24-Jun-2022		1-Jul-2022	NA	NA	Change 396.52

<a href="#">Monograph Title</a>	<a href="#">Section</a>	<a href="#">Source Publication</a>	<a href="#">Page Number</a>	<a href="#">Errata Post Date</a>	<a href="#">Sort ascending</a>	<a href="#">Errata Official Date</a>	<a href="#">Target Errata Print Publication</a>	<a href="#">Target Online Fix Publication</a>	Description
IVERMECTIN	CHEMICAL INFORMATION	USPNF Online	Online	24-Jun-2022		1-Jul-2022	NA	NA	to: 396.53 Change C <sub>48</sub> H <sub>74</sub> O <sub>14</sub> (Component H <sub>2</sub> B <sub>1a</sub> ) 875.09 C <sub>47</sub> H <sub>72</sub> O <sub>14</sub> (Component H <sub>2</sub> B <sub>1b</sub> ) 861.07 Component H <sub>2</sub> B <sub>1a</sub> : Avermectin A <sub>1a</sub> , 5-O -demethyl-22,23 -dihydro- (2aE,4E,8E )-(5?S,6S,6?R ,7S,11R,13R ,15S,17aR ,20R,20aR ,20bS )-6?-(S)-sec -Butyl-3?,4?,5?, 6,6?,7,10,11,14, 15,17a,20,20a,2 0b-tetradecahyd ro-20,20b-dihyd roxy[11,15-meth ano-2H,13H ,17H -furo[

<a href="#">Monograph Title</a> <a href="#">Section</a>	<a href="#">Source Publication</a>	<a href="#">Page Number</a>	<a href="#">Errata Post Date</a> <a href="#">Sort ascending</a>	<a href="#">Errata Official Date</a>	<a href="#">Target Errata Print Publication</a>	<a href="#">Target Online Fix Publication</a>	Description
							4,3,2- <i>pq</i> [[2,6]benzodiox acyclooctadecin -13,2?- [2 <i>H</i> ]pyran]-7-yl 2, 6-di deoxy- 4-O
							O -me thyl-?- L- <i>arabino</i> -hexopyranosyl) -3-O-methyl-?- L- <i>arabino</i> -hexopyranosid e CAS RN®: 70161-11-4. Component H <sub>2</sub> B <sub>1b</sub> : Avermectin A <sub>1a</sub> , 5-O -demethyl-25-d e(1-methylpropy l)-22,23-dihydro -25-(1-methylet

<a href="#">Monograph Title</a> <a href="#">Section</a>	<a href="#">Source Publication</a>	<a href="#">Page Number</a>	<a href="#">Errata Post Date</a> <a href="#">Sort ascending</a>	<a href="#">Errata Official Date</a>	<a href="#">Target Errata Print Publication</a>	<a href="#">Target Online Fix Publication</a>	Description
							hyl)- (2aE,4E,8E )-(5?S,6S,6?R ,7S,11R,13R ,15S,17aR ,20R,20aR ,20bS )-3?,4?,5?,6,6?, 7,10,11,-oxospir o[11,15-methan o-2H,13H,17H -furo[ 4,3,2-pq ][2,6]benzodiox acyclooctadecin -13,2?[ 2H]pyran]-7-yl 2, 6-di deoxy- 4-O  O -me thyl-?- L- <i>arabino</i> -hexopyranosyl) -3-O-methyl-?- L-

<a href="#">Monograph Title</a> <a href="#">Section</a>	<a href="#">Source Publication</a>	<a href="#">Page Number</a>	<a href="#">Errata Post Date</a> <a href="#">Sort ascending</a>	<a href="#">Errata Official Date</a>	<a href="#">Target Errata Print Publication</a>	<a href="#">Target Online Fix Publication</a>	Description
							<p><i>arabino</i>  -hexopyranosid  e CAS RN®:  70209-81-3;  UNII:  0W28CYI3TU.  to:  C<sub>48</sub>H<sub>74</sub>O<sub>14</sub>  (Component  H<sub>2</sub>B<sub>1a</sub>) 875.11  C<sub>47</sub>H<sub>72</sub>O<sub>14</sub>  (Component  H<sub>2</sub>B<sub>1b</sub>) 861.08  Component  H<sub>2</sub>B<sub>1a</sub>:  Avermectin A<sub>1a</sub>,  5-O  -demethyl-22,23  -dihydro-  (2a<i>E</i>,4<i>E</i>,8<i>E</i>  )-(5'<i>S</i>,6<i>S</i>,6'<i>R</i>  ,7<i>S</i>,11<i>R</i>,13<i>R</i>  ,15<i>S</i>,17a<i>R</i>  ,20<i>R</i>,20a<i>R</i>  ,20b<i>S</i>)-6'-(<i>S</i>)-  sec  -Butyl-3',4',5',  6,6',7,10,11,14,  15,17a,20,20a,2  0b-tetradecahyd  ro-20,20b-dihyd</p>



<a href="#">Monograph Title</a> <a href="#">Section</a>	<a href="#">Source Publication</a>	<a href="#">Page Number</a>	<a href="#">Errata Post Date</a> <a href="#">Sort ascending</a>	<a href="#">Errata Official Date</a>	<a href="#">Target Errata Print Publication</a>	<a href="#">Target Online Fix Publication</a>	Description
							roxy-5',6,8,19-t etramethyl-17-o xospiro [11,15- meth ano-2 <i>H</i> ,13 <i>H</i> ,17 <i>H</i> -furo[ 4,3,2- <i>pq</i> ][2,6]benzodiox acyclooctadecin -13,2'-[ 2 <i>H</i> ]pyran]-7-yl 2, 6-di deoxy- 4-O  O -me thyl-?- L- <i>arabino</i> -hexopyranosyl) -3-O-methyl-?- L- <i>arabino</i> -hexopyranosid e CAS RN®: 71827-03-7; UNII:

<a href="#">Monograph Title</a> <a href="#">Section</a>	<a href="#">Source Publication</a>	<a href="#">Page Number</a>	<a href="#">Errata Post Date</a> <a href="#">Sort ascending</a>	<a href="#">Errata Official Date</a>	<a href="#">Target Errata Print Publication</a>	<a href="#">Target Online Fix Publication</a>	Description
							91Y2202OUW. Component H <sub>2</sub> B <sub>1b</sub> : Avermectin A <sub>1a</sub> , 5-O -demethyl-25-d e(1-methylpropy l)-22,23-dihydro -25-(1-methylet hyl)- (2aE,4E,8E )-(5'S,6S,6'R ,7S,11R,13R ,15S,17aR ,20R,20aR ,20bS )-3',4',5',6,6', 7,10,11,14,15,1 7a,20,20a,20b- Tetradecahydro -20,20b-dihydro xy-6'-isopropyl- 5',6,8,19-tetra methyl-17-oxos piro [11,15-met hano-2H,13H ,17H -furo[ 4,3,2-pq ][2,6]benzodiox acyclooctadecin

<a href="#">Monograph Title</a> <a href="#">Section</a>	<a href="#">Source Publication</a>	<a href="#">Page Number</a>	<a href="#">Errata Post Date</a> <a href="#">Sort ascending</a>	<a href="#">Errata Official Date</a>	<a href="#">Target Errata Print Publication</a>	<a href="#">Target Online Fix Publication</a>	Description
DOXYCYCLINE ADDITIONAL RHYCLATE CAPSULES	USPNF Online Online <i>Reference Standards &lt;11&gt;</i>		24-Jun-2022	1-Jul-2022	NA	NA	-13,2'-[2H]pyran]-7-yl 2, 6-di deoxy-4-O O -me thyl-?- L- <i>arabino</i> -hexopyranosyl) -3-O-methyl-?- L- <i>arabino</i> -hexopyranosid e CAS RN®: 70209-81-3; UNII: 0W28CYI3TU. In USP Doxycycline Related Compound A RS: Change 444.43 to: 444.44 AND

<a href="#">Monograph Title</a> <a href="#">Section</a>	<a href="#">Source Publication</a>	<a href="#">Page Number</a>	<a href="#">Errata Post Date</a> <a href="#">Sort ascending</a>	<a href="#">Errata Official Date</a>	<a href="#">Target Errata Print Publication</a>	<a href="#">Target Online Fix Publication</a>	Description
							Change (4S,4aR,5S,5aR,6S,12aS)-4-(Dimethylamino)-1,4,4a,5,5a,6,11,12a-octahydro-3,5,10,12,12a-pentahydroxy-6-methyl-1,11-dioxo-2-naphthacenenecarboxamide, monohydrochloride. $C_{22}H_{24}N_2O_8 \cdot HCl$ 480.13 to: (4S,4aR,5S,5aR,6S,12aS)-4-(Dimethylamino)-1,4,4a,5,5a,6,11,12a-octahydro-3,5,10,12,12a-pentahydroxy-6-methyl-1,11-dioxo-2-naphthacenenecarboxamide hydrochloride. C

<a href="#">Monograph Title</a>	<a href="#">Section</a>	<a href="#">Source Publication</a>	<a href="#">Page Number</a>	<a href="#">Errata Post Date</a>	<a href="#">Sort ascending</a>	<a href="#">Errata Official Date</a>	<a href="#">Target Errata Print Publication</a>	<a href="#">Target Online Fix Publication</a>	Description
HYDROCODONE BITARTRATE AND HOMATROPINE METHYLBROMIDE TABLETS	ASSAY/ Procedure	USPNF Online	Online	24-Jun-2022		1-Jul-2022	NA	NA	$^{22}\text{H}_{24}\text{N}_2\text{O}_8 \cdot \text{HCl}$ 480.90 In Buffer: Change Adjust with phosphoric acid to a pH of $6.4 \pm 0.01$ . to: Adjust with phosphoric acid to a pH of $6.4 \pm 0.1$ .
NONOXYNOL 9	CHEMICAL INFORMATION	USPNF Online	Online	24-Jun-2022		1-Jul-2022	NA	NA	Update the chemical structure AND Change ?-(p -Nonylphenyl)-? -hydroxynona(oxyethylene) CAS RN®: 26027-38-3. to: ?-(4-Nonylphenyl)-?-hydroxynona(oxyethylene)
TRYPTOPHAN	ADDITIONAL REQUIREMENT	USPNF Online	Online	24-Jun-2022		1-Jul-2022	NA	NA	In USP Tryptophan

<a href="#">Monograph Title</a>	<a href="#">Section</a>	<a href="#">Source Publication</a>	<a href="#">Page Number</a>	<a href="#">Errata Post Date</a>	<a href="#">Sort ascending</a>	<a href="#">Errata Official Date</a>	<a href="#">Target Errata Print Publication</a>	<a href="#">Target Online Fix Publication</a>	Description
									<p>Related Compound A  RS: Change  3,3'-[Ethylidene bis(1<i>H</i>-indole-1,3-diyl)] bis[2<i>S</i>]-2-aminopropionic acid.  <math>C_{24}H_{26}N_4O_4</math>  432.49  to:  (2<i>S</i>,2'<i>S</i>)-3,3'-[Ethane-1,1-diyl bis(1<i>H</i>-indole-1,3-diyl)] bis(2-aminopropionic acid).  <math>C_{24}H_{26}N_4O_4</math>  434.50</p>
DOXYCYCLINE ADDITIONAL TABLETS	EQUIREMENT	USPNF Online	Online	24-Jun-2022		1-Jul-2022	NA	NA	<p>In USP  Doxycycline  Related Compound A  RS: Change  444.43  to:  444.44  AND  Change</p>
									<p><i>S/USP Reference Standards ?11?</i></p>
									<p><i>S/USP Reference Standards &lt;11&gt;</i></p>

<a href="#">Monograph Title</a> <a href="#">Section</a>	<a href="#">Source Publication</a>	<a href="#">Page Number</a>	<a href="#">Errata Post Date</a> <a href="#">Sort ascending</a>	<a href="#">Errata Official Date</a>	<a href="#">Target Errata Print Publication</a>	<a href="#">Target Online Fix Publication</a>	Description
							<p>(4<i>S</i>,4<i>aR</i>,5<i>S</i>,5<i>aR</i>,6<i>S</i>,12<i>aS</i>)-4-(Dimethylamino)-1,4,4<i>a</i>,5,5<i>a</i>,6,11,12<i>a</i>-octahydro-3,5,10,12,12<i>a</i>-pentahydroxy-6-methyl-1,11-dioxo-2-naphthacenenecarboxamide, monohydrochloride.  <math>C_{22}H_{24}N_2O_8 \cdot HCl</math> 480.13  to:  (4<i>S</i>,4<i>aR</i>,5<i>S</i>,5<i>aR</i>,6<i>S</i>,12<i>aS</i>)-4-(Dimethylamino)-1,4,4<i>a</i>,5,5<i>a</i>,6,11,12<i>a</i>-octahydro-3,5,10,12,12<i>a</i>-pentahydroxy-6-methyl-1,11-dioxo-2-naphthacenenecarboxamide hydrochloride.  <math>C_{22}H_{24}N_2O_8 \cdot HCl</math> 480.90</p>

<a href="#">Monograph Title</a>	<a href="#">Section</a>	<a href="#">Source Publication</a>	<a href="#">Page Number</a>	<a href="#">Errata Post Date</a>	<a href="#">Errata Official Date</a>	<a href="#">Target Errata Print Publication</a>	<a href="#">Target Online Fix Publication</a>	Description
MILK THISTLE CAPSULES	PERFORMANCE TESTS/ <i>Disintegration and Dissolution &lt;2040&gt;, Dissolution</i>	USPNF Online	Online	24-Jun-2022	1-Aug-2022	NA	NA	In <i>Medium</i> : Change <i>Buffer</i> containing 2% lauryl sulfate; 900 mL to: <i>Buffer</i> containing 2% sodium lauryl sulfate; 900 mL
AMMONIUM GLYCYRRHIZATE	ASSAY/ <i>Content of Ammonium 18?- and 18?- Glycyrrhizate</i>	USPNF Online	Online	24-Jun-2022	1-Jul-2022	NA	NA	In <i>Analysis</i> : Change $M_{W(Salt)}$ = molecular weight of ammonium glycyrrhizate, 840.08 g/mol $M_{W(Acid)}$ = molecular weight of glycyrrhizic acid, 821.59 g/mol to: $M_{W(Salt)}$ = molecular weight of ammonium glycyrrhizate,



<a href="#">Monograph Title</a> <a href="#">Section</a>	<a href="#">Source Publication</a>	<a href="#">Page Number</a>	<a href="#">Errata Post Date</a> <a href="#">Sort ascending</a>	<a href="#">Errata Official Date</a>	<a href="#">Target Errata Print Publication</a>	<a href="#">Target Online Fix Publication</a>	Description
MELENGESTROL ACETATE	<i>USP Reference standards ?11?</i> <i>USPNF Online</i>	Online	24-Jun-2022	1-Jul-2022	NA	NA	<p>839.97 g/mol  <math>M_{W(Acid)}</math> =  molecular weight of glycyrrhizic acid, 822.94 g/mol</p> <p>In USP Melengestrol Acetate Related Compound A  RS: Change 16-Methylene-17?-hydroxy-4-pregnene-3,20-dione 17-acetate.  to:  16-Methylene-3,20-dioxopregn-4-en-17-yl acetate.  AND  In USP Melengestrol Acetate Related Compound B  RS: Change 17?-Hydroxy-6,16-dimethylene progna-4-ene-3,20-dione</p>

<a href="#">Monograph Title</a> <a href="#">Section</a>	<a href="#">Source Publication</a>	<a href="#">Page Number</a>	<a href="#">Errata Post Date</a> <a href="#">Sort ascending</a>	<a href="#">Errata Official Date</a>	<a href="#">Target Errata Print Publication</a>	<a href="#">Target Online Fix Publication</a>	Description
DOXYCYCLINE ADDITIONAL REQUIREMENT S/USP Reference Standards <11>	USPNF Online	Online	24-Jun-2022	1-Jul-2022	NA	NA	17-acetate. to: 6,16-Dimethylene-3,20-dioxopregn-4-en-17-yl acetate. In USP Doxycycline Related Compound A RS: Change 444.43 to: 444.44 AND Change (4S,4aR,5S,5aR,6S,12aS)-4-(Dimethylamino)-1,4,4a,5,5a,6,11,12a-octahydro-3,5,10,12,12a-pentahydroxy-6-methyl-1,11-dioxo-2-naphthacene carboxamide monohydrochloride. C <sub>22</sub> H <sub>24</sub> N <sub>2</sub> O <sub>8</sub> · HCl 480.13

<a href="#">Monograph Title</a> <a href="#">Section</a>	<a href="#">Source Publication</a>	<a href="#">Page Number</a>	<a href="#">Errata Post Date</a> <a href="#">Sort ascending</a>	<a href="#">Errata Official Date</a>	<a href="#">Target Errata Print Publication</a>	<a href="#">Target Online Fix Publication</a>	Description
DOXYCYCLINE ADDITIONAL REQUIREMENT S/USP Reference Standards <11>	USPNF Online	Online	24-Jun-2022	1-Jul-2022	NA	NA	to: (4S,4aR,5S,5aR,6S,12aS)-4-(Dimethylamino)-1,4,4a,5,5a,6,11,12a-octahydro-3,5,10,12,12a-pentahydroxy-6-methyl-1,11-dioxo-2-naphthacene-carboxamide hydrochloride. C <sub>22</sub> H <sub>24</sub> N <sub>2</sub> O <sub>8</sub> · HCl 480.90 In USP Doxycycline Related Compound A RS: Change 444.43 to: 444.44 AND Change (4S,4aR,5S,5aR,6S,12aS)-4-(Dimethylamino)-1,4,4a,5,5a

<a href="#">Monograph Title</a>	<a href="#">Section</a>	<a href="#">Source Publication</a>	<a href="#">Page Number</a>	<a href="#">Errata Post Date</a>	<a href="#">Sort ascending</a>	<a href="#">Errata Official Date</a>	<a href="#">Target Errata Print Publication</a>	<a href="#">Target Online Fix Publication</a>	Description
HYDROCODONE BITARTRATE AND HOMATROPINE METHY	IM PURITIES/Limit of Homatropine Hydrobromide and Related	USPNF Online	Online	24-Jun-2022		1-Jul-2022	NA	NA	,6,11,12a-octahydro-3,5,10,12,12a-pentahydroxy-6-methyl-1,11-dioxo-2-naphthacenenecarboxamide, monohydrochloride. $C_{22}H_{24}N_2O_8 \cdot HCl$ 480.13 to: (4S,4aR,5S,5aR,6S,12aS)-4-(Dimethylamino)-1,4,4a,5,5a,6,11,12a-octahydro-3,5,10,12,12a-pentahydroxy-6-methyl-1,11-dioxo-2-naphthacenenecarboxamide hydrochloride. $C_{22}H_{24}N_2O_8 \cdot HCl$ 480.90 In Buffer: Change Adjust with phosphoric acid to a pH of 6.4 ±

<a href="#">Monograph Title</a>	<a href="#">Section</a>	<a href="#">Source Publication</a>	<a href="#">Page Number</a>	<a href="#">Errata Post Date</a>	<a href="#">Sort ascending</a>	<a href="#">Errata Official Date</a>	<a href="#">Target Errata Print Publication</a>	<a href="#">Target Online Fix Publication</a>	Description
LBROMIDE TABLETS	<i>Substances</i>								0.01. to: Adjust with phosphoric acid to a pH of 6.4 ± 0.1.
HOMATROPIN CHEMICAL E HYDROBRO MIDE	INFORMATION	<i>USPNF Online</i>	Online	24-Jun-2022		1-Jul-2022	NA	NA	Change 356.25 to: 356.26 AND Change 1?H,5?H-Tropan-3?-ol mandelate (ester) hydrobromide to: (1R,3r,5S)-8-Methyl-8-azabicyclo[3.2.1]octan-3-yl 2-hydroxy-2-phenylacetate hydrobromide
DOXYCYCLINE ADDITIONAL R HYCLATE DEL AYED- RELEASE TABLETS	EQUIREMENT <i>S/USP Reference Standards &lt;11&gt;</i>	<i>USPNF Online</i>	Online	24-Jun-2022		1-Jul-2022	NA	NA	In USP Doxycycline Related Compound A RS: Change 444.43

<a href="#">Monograph Title</a> <a href="#">Section</a>	<a href="#">Source Publication</a>	<a href="#">Page Number</a>	<a href="#">Errata Post Date</a> <a href="#">Sort ascending</a>	<a href="#">Errata Official Date</a>	<a href="#">Target Errata Print Publication</a>	<a href="#">Target Online Fix Publication</a>	Description
							<p>to:  444.44  AND  Change  (4S,4aR,5S  ,5aR,6S  ,12aS  )-4-(Dimethylam  ino)-1,4,4a,5,5a  ,6,11,12a-octah  ydro-3,5,10,12,  12a-pentahydro  xy-6-methyl-1,1  1-dioxo-2-napht  hacenicarboxa  mide, monohydr  ochloride.  C<sub>22</sub>H<sub>24</sub>N<sub>2</sub>O<sub>8</sub> ·  HCl 480.13</p> <p>to:  (4S,4aR,5S  ,5aR,6S  ,12aS  )-4-(Dimethylam  ino)-1,4,4a,5,5a  ,6,11,12a-octah  ydro-3,5,10,12,  12a-pentahydro  xy-6-methyl-1,1  1-dioxo-2-napht  hacenicarboxa</p>

<a href="#">Monograph Title</a>	<a href="#">Section</a>	<a href="#">Source Publication</a>	<a href="#">Page Number</a>	<a href="#">Errata Post Date</a>	<a href="#">Sort ascending</a>	<a href="#">Errata Official Date</a>	<a href="#">Target Errata Print Publication</a>	<a href="#">Target Online Fix Publication</a>	Description
MILK THISTLE TABLETS	PERFORMANCE TESTS/ <i>Disintegration and Dissolution &lt;2040&gt;, Dissolution</i>	USPNF Online	Online	24-Jun-2022		1-Aug-2022	NA	NA	mide hydrochloride. $C_{22}H_{24}N_2O_8 \cdot HCl$ 480.90 In <i>Medium</i> : Change <i>Buffer</i> containing 2% lauryl sulfate; 900 mL to: <i>Buffer</i> containing 2% sodium lauryl sulfate; 900 mL
AMMONIUM GLYCYRRHIZATE	SPECIFIC TESTS	USPNF Online	Online	24-Jun-2022		1-Jul-2022	NA	NA	In <i>Optical Rotation, Specific Rotation</i> ?781?: Change ?781? to: ?781S? AND Change <b>Acceptance criteria</b> : +49.0 to +55.0 on the anhydrous basis to:

<a href="#">Monograph Title</a> <a href="#">Section</a>	<a href="#">Source Publication</a>	<a href="#">Page Number</a>	<a href="#">Errata Post Date</a> <a href="#">Sort ascending</a>	<a href="#">Errata Official Date</a>	<a href="#">Target Errata Print Publication</a>	<a href="#">Target Online Fix Publication</a>	Description
METHOTREXATE ADDITIONAL REQUIREMENT S/USP Reference Standards ?11?	USPNF Online	Online	24-Jun-2022	1-Jul-2022	NA	NA	<b>Acceptance criteria:</b> +49.0° to +55.0° on the anhydrous basis In USP Methotrexate System Suitability Mixture RS: Change $C_{22}H_{26}N_8O_5 \cdot HCl$ 518.95 to: $C_{22}H_{26}N_8O_5 \cdot x HCl$
DOXYCYCLINE FOR INJECTION ADDITIONAL REQUIREMENT S/USP Reference Standards <11>	USPNF Online	Online	24-Jun-2022	1-Jul-2022	NA	NA	In USP Doxycycline Related Compound A RS: Change 444.43 to: 444.44 AND Change (4S,4aR,5S,5aR,6S,12aS)-4-(Dimethylamino)-1,4,4a,5,5a



<a href="#">Monograph Title</a>	<a href="#">Section</a>	<a href="#">Source Publication</a>	<a href="#">Page Number</a>	<a href="#">Errata Post Date</a>	<a href="#">Sort ascending</a>	<a href="#">Errata Official Date</a>	<a href="#">Target Errata Print Publication</a>	<a href="#">Target Online Fix Publication</a>	Description
ASCORBIC ACID INJECTION	ASSAY/ Procedure	USPNF Online	Online	24-Jun-2022		1-Jul-2022	NA	NA	,6,11,12a-octahydro-3,5,10,12,12a-pentahydroxy-6-methyl-1,11-dioxo-2-naphthacenenecarboxamide, monohydrochloride. C <sub>22</sub> H <sub>24</sub> N <sub>2</sub> O <sub>8</sub> · HCl 480.13 to: (4S,4aR,5S,5aR,6S,12aS)-4-(Dimethylamino)-1,4,4a,5,5a,6,11,12a-octahydro-3,5,10,12,12a-pentahydroxy-6-methyl-1,11-dioxo-2-naphthacenenecarboxamide hydrochloride. C <sub>22</sub> H <sub>24</sub> N <sub>2</sub> O <sub>8</sub> · HCl 480.90 In Chromatographic system/Column: Change

<a href="#">Monograph Title</a>	<a href="#">Section</a>	<a href="#">Source Publication</a>	<a href="#">Page Number</a>	<a href="#">Errata Post Date</a>	<a href="#">Sort ascending</a>	<a href="#">Errata Official Date</a>	<a href="#">Target Errata Print Publication</a>	<a href="#">Target Online Fix Publication</a>	Description
RIFABUTIN	CHEMICAL INFORMATION	USPNF Online	Online	27-May-2022		1-Jun-2022	NA	NA	150-cm x 6-mm; packing L39 to: 15-cm x 6-mm; packing L39 Change 847.00 to: 847.02 AND Change (9S,12E,14S,15R,16S,17R,18R,19R,20S,21S,22E,24Z)-6,16,18,20-Tetrahydroxy-1-isobutyl-14-methoxy-7,9,15,17,19,21,25-heptamethylspiro[9,4-(epoxypentadeca[1,11,13]trienimino)-2H-furo[2,3:7,8]naphth[1,2-d]imidazole-2,4-piperidine]-5,10,26-(3H,9H

<a href="#">Monograph Title</a> <a href="#">Section</a>	<a href="#">Source Publication</a>	<a href="#">Page Number</a>	<a href="#">Errata Post Date</a> <a href="#">Sort ascending</a>	<a href="#">Errata Official Date</a>	<a href="#">Target Errata Print Publication</a>	<a href="#">Target Online Fix Publication</a>	Description
							)-trione-16-acetate to: (9S,12E,14S,15R,16S,17R,18R,19R,20S,21S,22E,24Z)-6,18,20-Trihydroxy-1'-isobutyl-14-methoxy-7,9,15,17,19,21,25-heptamethyl-5,10,26-trioxo-3,5,9,10-tetrahydrospiro[9,4-(epoxypentadeca[1,11,13]trienimino)-2H-furo[2',3':7,8]naphthalen-1-ylidene-2-imidazole-2,4'-piperidin]-16-yl acetate Change 1025.87 to: 1025.88 In Acceptance criteria/Total impurities:
DOXYCYCLINE CHEMICAL HYCLATE INFORMATION	USPNF Online	Online	27-May-2022	1-Jun-2022	NA	NA	
FLUOCINOLON E ACETONIDE	Organic Impurities USPNF Online	Online	27-May-2022	1-Jun-2022	NA	NA	

<a href="#">Monograph Title</a> <a href="#">Section</a>	<a href="#">Source Publication</a>	<a href="#">Page Number</a>	<a href="#">Errata Post Date</a> <a href="#">Sort ascending</a>	<a href="#">Errata Official Date</a>	<a href="#">Target Errata Print Publication</a>	<a href="#">Target Online Fix Publication</a>	Description
NANDROLONE ASSAY/ DECANOATE <i>Procedure</i>	USPNF Online	Online	27-May-2022	1-Jun-2022	NA	NA	<p>Change NMT 2.5%. Disregard any peak below 0.05% of the peak area of fluocinolone acetonide from the <i>Standard solution</i>.</p> <p>to: NMT 2.5%. Disregard any peak below 0.05% of the peak area of fluocinolone acetonide from the <i>Sample solution</i>.</p> <p>In <i>Analysis</i>: Change Calculate the percentage of Nandrolone Decanoate (C<sub>28</sub>H<sub>44</sub>O<sub>3</sub>) in the portion of Nandrolone Decanoate taken:</p>

<a href="#">Monograph Title</a>	<a href="#">Section</a>	<a href="#">Source Publication</a>	<a href="#">Page Number</a>	<a href="#">Errata Post Date</a>	<a href="#">Sort ascending</a>	<a href="#">Errata Official Date</a>	<a href="#">Target Errata Print Publication</a>	<a href="#">Target Online Fix Publication</a>	Description
IBUPROFEN ORAL SUSPENSION	PERFORMANCE TESTS/ <i>Dissolution</i> <711>	USPNF Online	Online	27-May-2022		1-Jun-2022	NA	NA	<p>to:</p> <p>Calculate the percentage of nandrolone decanoate (C<sub>28</sub>H<sub>44</sub>O<sub>3</sub>) in the portion of Nandrolone Decanoate taken:</p> <p>In <i>Analysis</i>: Change Result = <math>(R_U/R_S) \times C_S \times V \times (D/W_U) \times (1/L) \times 100</math> <i>R<sub>U</sub></i> = peak area ratio of ibuprofen to benzophenone from the <i>Sample solution</i> <i>R<sub>S</sub></i> = peak area ratio of ibuprofen to benzophenone from the <i>Standard solution</i> <i>C<sub>S</sub></i> = concentration of</p>

<a href="#">Monograph Title</a> <a href="#">Section</a>	<a href="#">Source Publication</a>	<a href="#">Page Number</a>	<a href="#">Errata Post Date</a> <a href="#">Sort ascending</a>	<a href="#">Errata Official Date</a>	<a href="#">Target Errata Print Publication</a>	<a href="#">Target Online Fix Publication</a>	Description
							<p>USP Ibuprofen RS in the <i>Standard solution</i> (mg/mL)  <i>V</i> = volume of <i>Medium</i>, 900 mL  <i>D</i> = density of Oral Suspension (g/mL)  <i>W<sub>U</sub></i> = weight of the portion of Oral Suspension added to the <i>Medium</i> (g)  <i>L</i> = label claim (mg/mL)  to:  Result = <math>(R_U/R_S) \times C_S \times V \times (d/W_U) \times D \times (1/L) \times 100</math>  <i>R<sub>U</sub></i> = peak area ratio of ibuprofen to benzophenone from the <i>Sample solution</i></p>

<a href="#">Monograph Title</a> <a href="#">Section</a>	<a href="#">Source Publication</a>	<a href="#">Page Number</a>	<a href="#">Errata Post Date</a> <a href="#">Sort ascending</a>	<a href="#">Errata Official Date</a>	<a href="#">Target Errata Print Publication</a>	<a href="#">Target Online Fix Publication</a>	Description
							$R_S$ = peak area ratio of ibuprofen to benzophenone from the <i>Standard solution</i> $C_S$ = concentration of USP Ibuprofen RS in the <i>Standard solution</i> (mg/mL) $V$ = volume of <i>Medium</i> , 900 mL $d$ = density of Oral Suspension (g/mL) $W_U$ = weight of the portion of Oral Suspension added to the <i>Medium</i> (g) $D$ = dilution factor of the <i>Sample solution</i> , 2

<a href="#">Monograph Title</a>	<a href="#">Section</a>	<a href="#">Source Publication</a>	<a href="#">Page Number</a>	<a href="#">Errata Post Date</a>	<a href="#">Sort ascending</a>	<a href="#">Errata Official Date</a>	<a href="#">Target Errata Print Publication</a>	<a href="#">Target Online Fix Publication</a>	Description
APREPITANT CAPSULES	PERFORMANCE TESTS/ <i>Dissolution</i> <711>	USPNF Online	Online	27-May-2022		1-Jun-2022	NA	NA	L = label claim (mg/mL) Change <b>Test 1</b> Dilute 1 mL of phosphoric acid with water to 1 L. to: <b>Test 1</b> AND Change <b>Dilute phosphoric acid:</b> to: <b>Dilute phosphoric acid:</b> Dilute 1 mL of phosphoric acid with water to 1 L.
DOXYCYCLINE HYCLATE	IMPURITIES/ <i>Organic Impurities</i>	USPNF Online	Online	27-May-2022		1-Jun-2022	NA	NA	In <i>Table 2</i> , footnote b: Change (4S,4aR,5S,5aR,6R,12aS)-2-Acetyl-4-(di



<a href="#">Monograph Title</a> <a href="#">Section</a>	<a href="#">Source Publication</a>	<a href="#">Page Number</a>	<a href="#">Errata Post Date</a> <a href="#">Sort ascending</a>	<a href="#">Errata Official Date</a>	<a href="#">Target Errata Print Publication</a>	<a href="#">Target Online Fix Publication</a>	Description
							methylamino)-4 a,5a,6,12a-tetra hydro-3,5,10,12 ,12a-pentahydr oxy-6-methyl-tet racene-1,11-dio xo-2-naphthace necarboxamide. to: (4 <i>S</i> ,4 <i>aR</i> ,5 <i>S</i> ,5 <i>aR</i> ,6 <i>R</i> ,12 <i>aS</i> )-2-Acetyl-4-(di methylamino)-3, 5,10,12,12a-pe ntahydroxy-6-m ethyl-4a,5a,6,12 a-tetrahydrotetr ac ene- 1,11(4 <i>H</i> ,5 <i>H</i> )-dione.

## Pagination

- [First page](#) « [First](#)
- [Previous page](#) ‹ [Previous](#)
- ...
- [Page 2](#)
- [Page 3](#)
- [Page 4](#)

- 
- [Page 5](#)
  - [Page 6](#)
  - [Page 7](#)
  - [Page 8](#)
  - [Page 9](#)
  - [Page 10](#)
  - ...
  - [Next page Next ›](#)
  - [Last page Last »](#)