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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 Aminosalicyclic Acid Tablets will result in anything that contains “Aminosalicyclic” OR “Acid” OR “Tablets”
 - A search for “Aminosalicyclic Acid Tablets” will result in anything that specifically contains “Aminosalicyclic 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.

Monograph Title	Section	Source	Page Number	Errata Post	Errata Official	Target Errata	Target Online	Description
		Publication		Date	Date	Print Publication	Fix Publication	
ECONAZOLE	ADDITIONAL R	<i>USPNF Online</i>	Online	26-Aug-2022	1-Sep-2022	NA	<i>USPNF 2023</i>	In USP

Monograph Title	Section	Source Publication	Page Number	Errata Post Date	Errata Official Date	Target Errata Print Publication	Target Online Fix Publication	Description
NITRATE	EQUIREMENT S/USP Reference Standards <11>						Issue 2	Econazole Related Compound C RS: Change 1-(4-Chlorobenzyl)-3-{2-[4-chlorobenzyl)oxy]-2-(2,4-dichlorophenyl)ethyl}-1H-imidazol-3-ium nitrate (salt). C ₂₅ H ₂₁ Cl ₁₄ N ₃ O ₄ 569.26 to: 1-(4-Chlorobenzyl)-3-{2-[4-chlorobenzyl)oxy]-2-(2,4-dichlorophenyl)ethyl}-1H-imidazol-3-ium chloride. C ₂₅ H ₂₁ Cl ₅ N ₂ O 542.71
CHROMATOGRAPHY	ADJUSTMENT OF CHROMATOGRAPHIC CONDITIONS	USPNF Online	Online	26-Aug-2022	1-Dec-2022	NA	NA	In <i>Liquid Chromatography: Isocratic Elution/Injection volume:</i> Change

Monograph Title Section	Source Publication	Page Number	Errata Post Date Sort ascending	Errata Official Date	Target Errata Print Publication	Target Online Fix Publication	Description
PROMETHAZINE HYDROCHLORIDE IM PURITIES/ <i>Organic Impurities</i>	USPNF Online	Online	26-Aug-2022	1-Sep-2022	NA	NA	<p>L_2 = internal diameter of the column used (mm)</p> <p>dc_1 = particle size indicated in the monograph (μm)</p> <p>dc_2 = particle size of the column used (μm)</p> <p>to:</p> <p>L_2 = new column length (mm)</p> <p>dc_1 = column internal diameter indicated in the monograph (mm)</p> <p>dc_2 = new column internal diameter (mm)</p> <p>Change System suitability solution: 5 $\mu\text{g/mL}$ each of USP</p>

Monograph Title Section	Source Publication	Page Number	Errata Post Date Sort ascending	Errata Official Date	Target Errata Print Publication	Target Online Fix Publication	Description
							<p>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>Standard solution: 5 µg/mL of USP Promethazine Hydrochloride RS from the <i>Standard stock solution</i></p> <p>Sensitivity solution: 0.25 µg/mL of USP Promethazine Hydrochloride RS from the <i>Standard solution</i></p> <p>to:</p>

Monograph Title Section	Source Publication	Page Number	Errata Post Date Sort ascending	Errata Official Date	Target Errata Print Publication	Target Online Fix Publication	Description
							<p>System suitability solution: 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>Standard solution: 5 µg/mL of USP Promethazine Hydrochloride RS from the <i>Standard stock solution</i> in <i>Diluent</i></p> <p>Sensitivity solution: 0.25</p>

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METHYLENE BLUE	IMPURITIES/ <i>Organic Impurities</i>	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 solution in Diluent</i></p> <p>In footnote a in <i>Table 2</i>: Change 3-(Dimethylamino)-7-(methylamino)-phenothiazine-5-ium chloride. to: 3-(Dimethylamino)-7-(methylamino)phenothiazine-5-ium chloride.</p>
MELOXICAM ORAL SUSPENSION	PERFORMANCE TESTS/ <i>Dissolution <711></i>	USPNF Online	Online	29-Jul-2022	1-Aug-2022	NA	NA	<p>In <i>Analysis</i>: Change W_U = weight of the Oral Suspension taken (mg) to: W_U = weight of the Oral Suspension</p>

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LATANOPROS ASSAY/ T Procedure	USPNF Online	Online	29-Jul-2022	1-Dec-2022	NA	NA	<p>taken (g) Change Standard solution: 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>Sample solution: Transfer 2.0 mg/mL of Latanoprost into a suitable volumetric flask, dissolve in dehydrated alcohol equivalent to</p>

Monograph Title Section	Source Publication	Page Number	Errata Post Date Sort ascending	Errata Official Date	Target Errata Print Publication	Target Online Fix Publication	Description
							<p>20% of the final volume, and dilute with chromatographic hexane to volume.</p> <p>to:</p> <p>Standard solution: 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>Sample solution: 2.0 mg/mL of</p>

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FORSKOHLII	COMPOSITION <i>/Content of Forskolin</i>	USPNF Online	Online	29-Jul-2022		1-Aug-2022	NA	NA	<p>Latanoprost prepared as follows.</p> <p>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, 100 Å; packing L1</p>

Monograph Title	Section	Source Publication	Page Number	Errata Post Date	Sort ascending	Errata Official Date	Target Errata Print Publication	Target Online Fix Publication	Description
SUTURES--NE PROCEDURE EDLE ATTACHMENT		USPNF Online	Online	29-Jul-2022		1-Aug-2022	NA	NA	In Removable Needle Attachment: Change For USP sizes 5-0 through 2-0, to: For USP sizes 5-0 through 2,
BACLOFEN INJECTION	SPECIFIC TESTS	USPNF Online	Online	29-Jul-2022		1-Aug-2022	NA	NA	Change • Osmolality and Osmolarity ?785?, Osmolality: 270–320 mOsm/kg to: • Osmolality and Osmolarity ?785? Osmolality: 270–320 mOsm/kg
POWDERED FORSKOHLII	COMPOSITION /Content of Forskolin	USPNF Online	Online	29-Jul-2022		1-Aug-2022	NA	NA	In Chromatographic system: Change Column:

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POWDERED FORSKOHLII EXTRACT	COMPOSITION /Content of Forskolin	USPNF Online	Online	29-Jul-2022		1-Aug-2022	NA	NA	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 to: 396.53

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IVERMECTIN	CHEMICAL INFORMATION	USPNF Online	Online	24-Jun-2022		1-Jul-2022	NA	NA	Change $C_{48}H_{74}O_{14}$ (Component H_2B_{1a}) 875.09 $C_{47}H_{72}O_{14}$ (Component H_2B_{1b}) 861.07 Component H_2B_{1a} : Avermectin A _{1a} , 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[4,3,2-pq][2,6]benzodiox

Monograph Title Section	Source Publication	Page Number	Errata Post Date Sort ascending	Errata Official Date	Target Errata Print Publication	Target Online Fix Publication	Description
							<p>acyclooctadecin -13,2?- [2H]pyran]-7-yl 2, 6-di deoxy- 4-O</p> <p>O -me thyl-?- L- <i>arabino</i> -hexopyranosyl) -3-O-methyl-?- L- <i>arabino</i> -hexopyranosid e CAS RN®: 70161-11-4. Component H₂B_{1b}: Avermectin A_{1a}, 5-O -demethyl-25-d e(1-methylpropy l)-22,23-dihydro -25-(1-methylet hyl)-. (2aE,4E,8E</p>

Monograph Title Section	Source Publication	Page Number	Errata Post Date Sort ascending	Errata Official Date	Target Errata Print Publication	Target Online Fix Publication	Description
							<p>)-(5?S,6S,6?R,7S,11R,13R,15S,17aR,20R,20aR,20bS)</p> <p>-3?,4?,5?,6,6?,7,10,11,-oxospir</p> <p>o[11,15-methano-2H,13H,17H-furo[4,3,2-pq][2,6]benzodioxacyclooctadecin-13,2?[2H]pyran]-7-yl</p> <p>2,6-di</p> <p>deoxy-4-O</p> <p>O</p> <p>-me</p> <p>thyl-?-L-</p> <p><i>arabino</i></p> <p>-hexopyranosyl)</p> <p>-3-O-methyl-?-L-</p> <p><i>arabino</i></p> <p>-hexopyranosid</p>

Monograph Title Section	Source Publication	Page Number	Errata Post Date Sort ascending	Errata Official Date	Target Errata Print Publication	Target Online Fix Publication	Description
							<p>e CAS RN®: 70209-81-3; UNII: 0W28CYI3TU. to: C₄₈H₇₄O₁₄ (Component H₂B_{1a}) 875.11 C₄₇H₇₂O₁₄ (Component H₂B_{1b}) 861.08 Component H₂B_{1a}: Avermectin A_{1a}, 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-5',6,8,19-t etramethyl-17-o</p>

Monograph Title Section	Source Publication	Page Number	Errata Post Date Sort ascending	Errata Official Date	Target Errata Print Publication	Target Online Fix Publication	Description
							<p>xospiro [11,15-methano-2<i>H</i>,13<i>H</i>,17<i>H</i>]-furo[4,3,2-<i>pq</i>][2,6]benzodioxacyclooctadecin-13,2'-[2<i>H</i>]pyran]-7-yl 2,6-di-deoxy-4-O</p> <p>O</p> <p>-methyl-?-L-<i>arabino</i>-hexopyranosyl)-3-O-methyl-?-L-<i>arabino</i>-hexopyranoside CAS RN®: 71827-03-7; UNII: 91Y2202OUW. Component</p>

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							<p>H₂B_{1b}: Avermectin A_{1a}, 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 -13,2'-[2H]pyran]-7-yl</p>

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DOXYCYCLINE ADDITIONAL RHYCLATE EQUIREMENT CAPSULES S/USP Reference Standards <11>	USPNF Online	Online	24-Jun-2022	1-Jul-2022	NA	NA	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 Change (4S,4aR,5S

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HYDROCODO ASSAY/	USPNF Online	Online	24-Jun-2022	1-Jul-2022	NA	NA	,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 ₂₂ H ₂₄ N ₂ O ₈ · 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-naphthacene-carboxamide hydrochloride. C ₂₂ H ₂₄ N ₂ O ₈ · HCl 480.90 In Buffer.

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NE BITARTRATE AND HOMATR OPINE METHY LBROMIDE TABLETS	<i>Procedure</i>								Change Adjust with phosphoric acid to a pH of 6.4 ± 0.01. to: Adjust with phosphoric acid to a pH of 6.4 ± 0.1.
NONOXYNOL 9 CHEMICAL INFORMATION		<i>USPNF Online</i>	Online	24-Jun-2022		1-Jul-2022	NA	NA	Update the chemical structure AND Change ?-(p -Nonylphenyl)-? -hydroxynona(o xyethylene) CAS RN@: 26027-38-3. to: ?-(4-Nonylphen yl)-?-hydroxyno na(oxyethylene) .
TRYPTOPHAN	ADDITIONAL R EQUIREMENT S/USP <i>Reference</i> <i>Standards ?11?</i>	<i>USPNF Online</i>	Online	24-Jun-2022		1-Jul-2022	NA	NA	In USP Tryptophan Related Compound A RS: Change

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DOXYCYCLINE ADDITIONAL R TABLETS	<i>USPNF Online</i> Online		24-Jun-2022	1-Jul-2022	NA	NA	<p>3,3'-[Ethylidene bis(1<i>H</i>-indole-1,3-diyl)] bis[2<i>S</i>]-2-aminopropionic acid. $C_{24}H_{26}N_4O_4$ 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). $C_{24}H_{26}N_4O_4$ 434.50 In USP Doxycycline Related Compound A RS: Change 444.43 to: 444.44 AND Change (4<i>S</i>,4<i>aR</i>,5<i>S</i>,5<i>aR</i>,6<i>S</i>,12<i>aS</i>)</p>

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MILK THISTLE CAPSULES	PERFORMANCE TESTS/	USPNF Online	Online	24-Jun-2022		1-Aug-2022	NA	NA	<p>)-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</p> <p>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</p> <p>In <i>Medium</i>: Change <i>Buffer</i></p>

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									containing 2% lauryl sulfate; 900 mL to: <i>Buffer</i> containing 2% sodium lauryl sulfate; 900 mL
AMMONIUM G LYCYRRHIZATE	ASSAY/Content of Ammonium 18?- and 18?- Glycyrrhizate	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, 839.97 g/mol $M_{W(Acid)}$ = molecular

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MELENGESTR USP Reference OL ACETATE standards ?11?	USPNF Online	Online	24-Jun-2022	1-Jul-2022	NA	NA	weight of glycyrrhizic acid, 822.94 g/mol 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 17-acetate. to: 6,16-Dimethyle

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DOXYCYCLINE ADDITIONAL REQUIREMENTS/USP Reference Standards <11>	USPNF Online	Online	24-Jun-2022	1-Jul-2022	NA	NA	ne-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 ₂₂ H ₂₄ N ₂ O ₈ · HCl 480.13 to: (4S,4aR,5S,5aR,6S

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DOXYCYCLINE ADDITIONAL REQUIREMENT S/USP Reference Standards <11>	USPNF Online	Online	24-Jun-2022	1-Jul-2022	NA	NA	,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 ₂₂ H ₂₄ N ₂ O ₈ · 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,6,11,12a-octahydro-3,5,10,12,12a-pentahydro

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HYDROCODONE BITARTRATE AND HOMATROPINE METHYLBROMIDE TABLETS	IM PURITIES/Limit of Homatropine Hydrobromide and Related Substances	USPNF Online	24-Jun-2022	1-Jul-2022	NA	NA	xy-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 ± 0.01. to: Adjust with

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HOMATROPIN CHEMICAL E HYDROBRO INFORMATION MIDE	<i>USPNF Online</i> Online		24-Jun-2022	1-Jul-2022	NA	NA	phosphoric acid to a pH of 6.4 ± 0.1. Change 356.25 to: 356.26 AND Change 1?H,5?H- Tropan-3?-ol mandelate (ester) hydrobromide to: (1 <i>R</i> ,3 <i>r</i> ,5 <i>S</i>)-8-Methyl-8-az abicyclo[3.2.1]o ctan-3-yl 2-hydr oxy-2-phenylac etate hydrobromide In USP Doxycycline Related Compound A RS: Change 444.43 to: 444.44 AND
DOXYCYCLINE ADDITIONAL R HYCLATE DEL EQUIREMENT AYED- RELEASE TABLETS	<i>USPNF Online</i> Online <i>S/USP</i> <i>Reference</i> <i>Standards <11></i>		24-Jun-2022	1-Jul-2022	NA	NA	

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							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

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MILK THISTLE TABLETS	PERFORMANCE TESTS/ <i>Disintegration and Dissolution <2040>, Dissolution</i>	USPNF Online	Online	24-Jun-2022		1-Aug-2022	NA	NA	$^{22}\text{H}_{24}\text{N}_2\text{O}_8 \cdot \text{HCl}$ 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 In <i>Optical Rotation, Specific Rotation ?781?</i> : Change ?781? to: ?781S? AND Change Acceptance criteria : +49.0 to +55.0 on the anhydrous basis to: Acceptance criteria : +49.0°
AMMONIUM G LYCYRRHIZATE	SPECIFIC TESTS	USPNF Online	Online	24-Jun-2022		1-Jul-2022	NA	NA	

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METHOTREXATE	ADDITIONAL REQUIREMENT S/USP Reference Standards ?11?	USPNF Online	Online	24-Jun-2022	1-Jul-2022	NA	NA	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,6,11,12a-octahydro-3,5,10,12,

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ASCORBIC ACID INJECTION	ASSAY/ Procedure	USPNF Online	Online	24-Jun-2022		1-Jul-2022	NA	NA	<p>12a-pentahydroxy-6-methyl-1,11-dioxo-2-naphthacenenecarboxamide, monohydrochloride. $C_{22}H_{24}N_2O_8 \cdot HCl$ 480.13</p> <p>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</p> <p>In <i>Chromatographic system/Column: Change 150-cm x 6-mm; packing</i></p>

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RIFABUTIN	CHEMICAL INFORMATION	USP <i>NF Online</i>	Online	27-May-2022		1-Jun-2022	NA	NA	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)-trione-16-acetate

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							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]naphtho[1,2-d]imidazole-2,4'-piperidin]-16-yl acetate
DOXYCYCLINE CHEMICAL HYCLATE INFORMATION	USPNF Online	Online	27-May-2022	1-Jun-2022	NA	NA	Change 1025.87
FLUOCINOLON E ACETONIDE	Organic Impurities USPNF Online	Online	27-May-2022	1-Jun-2022	NA	NA	to: 1025.88 In Acceptance criteria/Total impurities: Change NMT 2.5%.

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NANDROLONE ASSAY/ DECANOATE <i>Procedure</i>	<i>USPNF Online</i> Online		27-May-2022	1-Jun-2022	NA	NA	<p>Disregard any peak below 0.05% of the peak area of fluocinolone acetonide from the <i>Standard solution</i>. to: NMT 2.5%.</p> <p>Disregard any peak below 0.05% of the peak area of fluocinolone acetonide from the <i>Sample solution</i>. In <i>Analysis</i>: Change Calculate the percentage of Nandrolone Decanoate (C₂₈H₄₄O₃) in the portion of Nandrolone Decanoate taken: to: Calculate the</p>

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IBUPROFEN ORAL SUSPENSION	PERFORMANCE TESTS/ <i>Dissolution</i> <711>	USPNF Online	Online	27-May-2022		1-Jun-2022	NA	NA	percentage of nandrolone decanoate (C ₂₈ H ₄₄ O ₃) in the portion of Nandrolone Decanoate taken: In <i>Analysis</i> : Change Result = $(R_U/R_S) \times C_S \times V \times (D/W_U) \times (1/L) \times 100$ <i>R_U</i> = peak area ratio of ibuprofen to benzophenone from the <i>Sample solution</i> <i>R_S</i> = peak area ratio of ibuprofen to benzophenone from the <i>Standard solution</i> <i>C_S</i> = concentration of USP Ibuprofen RS in the

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							<p><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_U</i> = weight of the portion of Oral Suspension added to the <i>Medium</i> (g) <i>L</i> = label claim (mg/mL) to: Result = $(R_U/R_S) \times C_S \times V \times (d/W_U) \times D \times (1/L) \times 100$ <i>R_U</i> = peak area ratio of ibuprofen to benzophenone from the <i>Sample solution</i> <i>R_S</i> = peak area ratio of</p>

Monograph Title Section	Source Publication	Page Number	Errata Post Date Sort ascending	Errata Official Date	Target Errata Print Publication	Target Online Fix Publication	Description
							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 $L =$ label claim (mg/mL)

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APREPITANT CAPSULES	PERFORMANCE TESTS/ <i>Dissolution</i> <711>	USPNF Online	Online	27-May-2022	1-Jun-2022	NA	NA	Change Test 1 Dilute 1 mL of phosphoric acid with water to 1 L. to: Test 1 AND Change Dilute phosphoric acid: to: Dilute phosphoric acid: 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 (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)-4 <i>a</i> ,5 <i>a</i> ,6,12 <i>a</i> -tetra

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DIMENHYDRIN OTHER COMP ATE TABLETS ONE NTS/ 8-Chlorotheoph ylline	USPNF Online	Online	27-May-2022	1-Jun-2022	NA	NA	hydro-3,5,10,12,12a-pentahydroxy-6-methyl-tetrahydro-1,11-dioxo-2-naphthacecarboxamide. to: (4S,4aR,5S,5aR,6R,12aS)-2-Acetyl-4-(dimethylamino)-3,5,10,12,12a-pentahydroxy-6-methyl-4a,5a,6,12a-tetrahydrotracene-1,11(4H,5H)-dione. In <i>Analysis</i> : Change C_U = nominal concentration of dimenhydrinate in the <i>Sample solution</i> (mg/mL) to: C_U = determined

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AZITHROMYCIN FOR ORAL SUSPENSION	ADDITIONAL REQUIREMENT S/USP Reference Standards <11>	USPNF Online	Online	27-May-2022		1-Jun-2022	NA	NA	concentration of dimenhydrinate in the <i>Sample solution</i> , as obtained in the Assay (mg/mL) In USP Azithromycin Related Compound F RS: Change 762.97 to: 762.98 AND In USP Desomethylazithromycin in RS: Change 590.79 to: 590.80
ATRACURIUM BESYLATE	CHEMICAL INFORMATION	USPNF Online	Online	27-May-2022		1-Jun-2022	NA	NA	Change 2-(2-Carboxyethyl)-1,2,3,4-tetrahydro-6,7-dimethoxy-2-methyl-1-veratrylisoquinolinium benzenesulfonate, pentamethylene ester

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CARBAMAZEPINE TABLETS ADDITIONAL REQUIREMENTS S/USP Reference Standards <11>	USP NF Online	Online	27-May-2022	1-Jun-2022	NA	NA	to: 2-(2-Carboxethyl)-1,2,3,4-tetrahydro-6,7-dimethoxy-2-methyl-1-veratrylisoquinolinium benzenesulfonate, pentamethylene ester In USP Carbamazepine Related Compound A RS: Change 238.28 to: 238.29 AND In USP Carbamazepine Related Compound B RS: Change 193.24 to: 193.25 AND In USP 9-Methylacridine RS: Change

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BOTANICAL EXTRACTS	PREPARATION S	USPNF Online	Online	27-May-2022		1-Jun-2022	NA	NA	<p>193.24 to: 193.25</p> <p>In <i>General Pharmacopeial R equ irement s/Pesticide Residues: Change where L is the limit in the original article as listed in Table 4 (see Pesticide Residue Analysis under Articles of Botanical Origin <561>)</i></p> <p>to: where L is the limit in the original article as listed in Table 5 (see Pesticide Residue Analysis under</p>

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DOXYCYCLINE ADDITIONAL REQUIREMENT S/USP Reference Standards <11>	USPNF Online	Online	27-May-2022	1-Jun-2022	NA	NA	Articles of Botanical Origin <561>) In USP Doxycycline Related Compound A RS: Change 444.43 to: 444.44 AND Change 480.13 to: 480.90	
MUPIROCIN NASAL OINTMENT	Related com pounds/	USPNF Online	Online	29-Apr-2022	1-May-2022	NA	NA	In footnote 2: Change 9-((E)-4-[(2R,3aS,6S,7S,8aRS)-2-((1RS,2S,3S)-1,3-Dihydroxy-2-methylbutyl)-7-hydroxyhexahydro-2H-furo[3,2-c]pyran-6-yl]-3-methylbut-2-enoyl

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									oxy}nonanoic acid. to: 9-{{E)-4-[(2R,3aS,6S,7S)-2-{{(1RS,2S,3S)-1,3-Dihydroxy-2-methylbutyl}-7-hydroxyhexahydro-2H-furo[3,2-c]pyran-6-yl]-3-methylbut-2-enoyloxy}nonanoic acid.
ACYCLOVIR	ASSAY/ Procedure	USPNF Online	Online	29-Apr-2022		1-May-2023	NA	NA	In the <i>Sample solution</i> : Change 0.1 N sodium hydroxide to: 0.01 N sodium hydroxide
DACARBAZINE IMPURITIES FOR INJECTION		USPNF Online	Online	29-Apr-2022		1-May-2022	NA	NA	Delete <i>Limit of 2-Azahypoxanthine</i> test
MUPIROCIN	IM	USPNF Online	Online	29-Apr-2022		1-May-2022	NA	NA	In footnote b:

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OINTMENT	PUR	ITIES/ <i>Organic Impurities/ Table</i>	2						Change 9- <i>(E)</i>)-4-[(2 <i>R</i> ,3 <i>aS</i> ,6 <i>S</i> ,7 <i>S</i> ,8 <i>aRS</i>)-2- <i>(1RS,2S</i> , <i>3S</i>)-1,3-Dihydroxy- 2-methylbutyl]-7 -hydroxyhexahy dro-2 <i>H</i> -furo[3, 2- <i>c</i>]pyran-6-yl]-3-m ethylbut-2-enoyl oxy}nonanoic acid. to: 9- <i>(E)</i>)-4-[(2 <i>R</i> ,3 <i>aS</i> ,6 <i>S</i> ,7 <i>S</i>)-2- <i>(1RS,2S</i> , <i>3S</i>)-1,3-Dihydroxy- 2-methylbutyl]-7 -hydroxyhexahy dro-2 <i>H</i> -furo[3, 2- <i>c</i>]pyran-6-yl]-3-m ethylbut-2-enoyl

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MECAMYLAMINE HYDROCHLORIDE	USP Reference Standards <11>	USPNF Online	29-Apr-2022	1-May-2022	NA	NA	oxy}nonanoic acid. In USP Mecamylamine Related Compound A RS: Change <i>N</i> ,1,7,7-Tetramethyl bicyclo [2.2.1]heptan-2-amine. C ₁₁ H ₂₁ N 167.29 to: <i>N</i> ,1,7,7-Tetramethyl bicyclo[2.2.1]heptan-2-amine hydrochloride. C ₁₁ H ₂₁ N · HCl 203.75

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