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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.
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- **Downloading:** You can download the Errata table in Comma-separated Value (.csv). The download will include the Errata that you have filtered on.
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Monograph Title	Section	Source	Page Number	Errata Post Date	Errata Official Date	Target Errata Print Publication	Target Online Fix Publication	Description
ORPHENADRI	ADDITIONAL R	USPNF Online	Online	28-Jan-2022	1-Feb-2022	NA	NA	In USP

Monograph Title	Section	Source Publication	Page Number	Errata Post Date	Sort ascending	Errata Official Date	Target Errata Print Publication	Target Online Fix Publication	Description
NE CITRATE INJECTION	EQUIREMENT <i>S/USP Reference Standards</i> <11>								Orphenadrine Related Compound B RS: Change <i>N</i> -Ethyl- <i>N,N</i> -dimethyl [2-(2-methylbenzhydryloxy)ethyl]ammonium chloride; also known as <i>N</i> -ethyl- <i>N,N</i> -dimethyl-2-[phenyl(o-tolyl)methoxy]ethanaminium chloride. to: <i>N</i> -Ethyl- <i>N,N</i> -dimethyl [2-(2-methylbenzhydryloxy)ethyl]ammonium chloride; also known as <i>N</i> -Ethyl- <i>N,N</i> -dimethyl-2-[phenyl(2-tolyl)methoxy]ethanaminium chloride.
NICARDIPINE HYDROCHLORIDE INJECTION	DEFINITION	<i>USPNF Online</i> Online		28-Jan-2022		1-Feb-2022	NA	NA	Change NTL 90.0% to:

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AMLODIPINE BESYLATE	Assay	USPNF Online	Online	28-Jan-2022	1-Feb-2022	NA	NA	NLT 90.0% In <i>Chromatographic system:</i> Change Chromatograph the <i>Standard preparation</i> , and record the peak responses as directed for <i>Procedure:</i> the standard deviation for replicate injections is not more than 2.0%. to: Chromatograph the <i>Standard preparation</i> , and record the peak responses as directed for <i>Procedure:</i> the relative standard deviation for replicate injections is not

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ONDANSETRO USP Reference N INJECTION standards <11>	USPNF Online	Online	28-Jan-2022	1-Feb-2022	NA	NA	more than 2.0%. In USP Ondansetron Related Compound A RS: Change 3-[(Dimethylamino)methyl]-1,2,3,9-tetrahydro-9-methyl-4H-carbazol-4-one hydrochloride. to: 3-[(Dimethylamino)methyl]-9-methyl-1,2,3,9-tetrahydro-4H-carbazol-4-one hydrochloride. AND In USP Ondansetron Related Compound C RS: Change 1,2,3,9-Tetrahydro-

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							9-me thyl-4 <i>H</i> -carbazol-4-one . to: 9-Methyl-1,2,3,9 -tet rahydr o-4 <i>H</i> -carbazol-4-one . AND In USP Ondansetron Related Compound D RS: Change 1,2,3,9-Tetrahy dro-9-methyl-3- met hylene- 4 <i>H</i> -carbazol-4-one . to: 9-Methyl-3-met hylene-1,2,3,9-t etra hydro-4 <i>H</i> -carbazol-4-one .

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LORATADINE CAPSULES	PERFORMANCE TESTS/ Dissolution <711>	USPNF Online	Online	28-Jan-2022		1-Feb-2022	NA	NA	In Tier 1/Solution A and Tier 2/Solution B: Change Tween 20 to: polysorbate 20
FULVESTRANT	CHEMICAL INFORMATION	USPNF Online	Online	28-Jan-2022		1-Feb-2022	NA	NA	Change 606.77 to: 606.78
ORPHENADRINE CITRATE EXTENDED-RELEASE TABLETS	ADDITIONAL REQUIREMENT S/USP Reference Standards <11>	USPNF Online	Online	28-Jan-2022		1-Feb-2022	NA	NA	In USP Orphenadrine Related Compound B RS: Change N-Ethyl-N,N-dimethyl [2-(methylbenzhydryl oxy)ethyl]ammonium chloride; Also known as N-Ethyl-N,N-dimethyl-2-[phenyl(o-tolyl)methoxy]ethanaminium chloride. to: N-Ethyl-N,N

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WARFARIN SODIUM TABLETS	ADDITIONAL REQUIREMENTS/USP Reference Standards <11>	USPNF Online	Online	28-Jan-2022		1-Feb-2022	NA	NA	-dimethyl [2-(2-methylbenzhydryloxy)ethyl]ammonium chloride; also known as N-Ethyl-N,N-dimethyl-2-[phenyl(2-tolyl)methoxy]ethanaminium chloride. In USP Warfarin Related Compound A RS: Change 3-(o-Hydroxyphenyl)-5-phenyl-2-cyclohexen-1-one. to: 3-(2-Hydroxyphenyl)-5-phenyl-2-cyclohexen-1-one. AND Change 264.33 to: 264.32 In Chromatographic system:
AMLODIPINE BESYLATE	Related compounds/Test 2	USPNF Online	Online	28-Jan-2022		1-Feb-2022	NA	NA	

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							Change Chromatograph the <i>Standard solution</i> , and record the peak responses as directed for <i>Procedure</i> : the standard deviation for replicate injections is not more than 10.0%. to: Chromatograph the <i>Standard solution</i> , and record the peak responses as directed for <i>Procedure</i> : the relative standard deviation for replicate injections is not more than 10.0%. Change USP
ONDANSETRO USP Reference N	USPNF 2021 Online standards <11> ISSUE 1		31-Dec-2021	1-Jan-2022	NA	NA	

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							Ondansetron RS USP Ondansetron Related Compound C RS 1,2,3,9-Tetrahy d ro- 9-me thyl-4 <i>H</i> -carbazol-4-one . USP Ondansetron Related Compound D RS 1,2,3,9-Tetrahy dro-9-methyl-3- met hylene- 4 <i>H</i> -carbazol-4-one . to: USP Ondansetron RS USP

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							Ondansetron Related Compound A RS 3-[(Dimethylamino)methyl]-9-methyl-1,2,3,9-tetrahydro-4H-carbazol-4-one hydrochloride. USP Ondansetron Related Compound C RS 9-Methyl-1,2,3,9-tetrahydro-4H-carbazol-4-one . USP Ondansetron Related Compound D RS 9-Methyl-3-methylene-1,2,3,9-tetrahydro-4H

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CARBOPLATIN IM FOR INJECTION	PURITIES/ <i>Limit of 1,1-Cyclobutanedicarboxylic Acid</i>	USPNF 2021 ISSUE 1	Online	31-Dec-2021		1-Jan-2022	NA	NA	-carbazol-4-one . In <i>Chromatographic system/Column:</i> Change 4.0-mm × 30-cm; packing L1 to: 3.9-mm × 30-cm; packing L1
CALCIUM SACCHARATE	ASSAY/ <i>Procedure</i>	USPNF 2021 ISSUE 1	Online	31-Dec-2021		1-Jan-2022	NA	NA	In <i>Analysis:</i> Change Result = $\{(V_S ? V_B) \times N \times F\} / W \times 100$ to: Result = $\{(V_S ? V_B) \times M \times F\} / W \times 100$ AND Change <i>N</i> = actual normality of the <i>Titrant</i> (mEq/mL) <i>F</i> = equivalency factor, 320.2

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EXTENDED PHENYTOIN SODIUM CAPSULES	ASSAY/ Procedure	<i>Revision Bulletin (Official March 01, 2021)</i>	Online	31-Dec-2021		1-Jan-2022	NA	NA	mg/mEq to: M = actual molarity of the <i>Titrant</i> (mmol/mL) F = equivalency factor, 320.26 mg/mmol In <i>Sample solution</i> : Change Nominally 0.6 mg/mL of phenytoin to: Nominally 0.6 mg/mL of phenytoin sodium AND In <i>Analysis</i> : Change C_U = nominal concentration of phenytoin in the <i>Sample solution</i> (mg/mL) to: C_U = nominal concentration of

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CARBOPLATIN IM		USPNF 2021	Online	31-Dec-2021		1-Jan-2022	NA	NA	phenytoin sodium in the <i>Sample solution</i> (mg/mL)
	PURITIES/Limit of 1,1-Cyclobutanedicarboxylic Acid	ISSUE 1							In <i>Chromatographic system/Column:</i> Change 4.0-mm × 30-cm; packing L1 to: 3.9-mm × 30-cm; packing L1
SUCCINYLCHOLINE CHLORIDE	ADDITIONAL REQUIREMENT S/USP Reference Standards <11>	USPNF 2021	Online	31-Dec-2021		1-Jan-2022	NA	NA	In USP Succinyl monocholine Chloride RS: Change Ethanaminium, 2-(carboxy-1-oxoproxy)-N,N,N-trimethyl-, chloride. to: Ethanaminium, 2-(3-carboxy-1-oxoproxy)-N,N,N-trimethyl-, chloride.

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ATORVASTATIN CALCIUM	ADDITIONAL REQUIREMENT S/USP Reference Standards <11>	USPNF 2021 ISSUE 1	Online	31-Dec-2021		1-Jan-2022	NA	NA	poxy)-N,N,N-trimethyl-, chloride; Also known as 2-[(3-Carboxypropyl)oxy]-N,N,N-trimethylethan-1-aminium chloride. In USP Atorvastatin Related Compound C RS: Change $C_{66}H_{66}F_4N_4O_{10}$ to: $C_{66}H_{66}CaF_4N_4O_{10}$ AND In USP Atorvastatin Related Compound E RS: Change 1155.38 to: 1155.36 In both Calculations in
EXTENDED PHENYTOIN	IMPUR	Revision Bulletin (Official	Online	31-Dec-2021		1-Jan-2022	NA	NA	

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SODIUM CAPSULES	ITIES/Organic Impurities	March 01, 2021)						Analysis: Change C_U = nominal concentration of phenytoin in the <i>Sample solution</i> ($\mu\text{g/mL}$) to: C_U = nominal concentration of phenytoin sodium in the <i>Sample solution</i> ($\mu\text{g/mL}$)
PRAZOSIN HYDROCHLORIDE	IMPURITIES/Organic Impurities	USPNF 2021 ISSUE 1	Online	31-Dec-2021	1-Jan-2022	NA	NA	In <i>Sensitivity solution</i> : Change 0.05 $\mu\text{g/mL}$ of USP Prazosin Hydrochloride RS in <i>Mobile phase</i> from the <i>Standard solution</i> to: 0.5 $\mu\text{g/mL}$ of USP Prazosin Hydrochloride RS in <i>Mobile phase</i> from the <i>Standard</i>

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CAPTOPRIL	ADDITIONAL REQUIREMENT S/USP Reference Standards <11>	USPNF 2021 ISSUE 1	Online	19-Nov-2021		1-Dec-2021	NA	NA	<i>solution</i> In USP Captopril Disulfide RS: Change L-Proline, 1,1?-[dithiobis(2-methyl-1-oxo-3,1-propanediyl)]bis-[S-(R*,R*)]-. to: (2?S)-[(2S,2?S)-3,3?-Disulfanediy]bis(2-methylpropanoyl)] di-L-proline.
ELEUTHERO ROOT AND RHIZOME DRY EXTRACT	ADDITIONAL REQUIREMENT S/USP Reference Standards <11>	USPNF 2021 ISSUE 1	Online	19-Nov-2021		1-Dec-2021	NA	NA	Change USP Eleutheroside B RS ?-D-Glucopyranoside, 4-(3-hydroxy-1-propenyl)-2,6-dimethoxyphenyl. C₁₇H₂₄O₉ 372.37 USP

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									Eleutheroside E RS ?-D-Glucopyranoside, (tetrahydr o-1 <i>H</i> ,3 <i>H</i> -furo(3,4-c)furan-1,4-diyl)bis(2,6-dimethoxy-4,1-pHENYLENE)bis-. $C_{34}H_{46}O_{18}$ 742.70 to: USP Eleutheroside B RS USP Eleutheroside E RS
DACARBAZINE IM FOR INJECTION	PURITIES/ <i>Organic Impurities</i>	USPNF 2021 ISSUE 1	Online	19-Nov-2021		1-Dec-2021	NA	NA	Delete <i>Diluent</i> : 4.0 mg/mL of citric acid in water
ELEUTHERO ROOT AND RHIZOME	ADDITIONAL REQUIREMENT S/USP Reference Standards <11>	USPNF 2021 ISSUE 1	Online	19-Nov-2021		1-Dec-2021	NA	NA	Change USP Eleutheroside B RS ?-D-Glucopyranoside, 4-(3-hydroxy-1-propenyl)

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							-2,6-dimethoxyp henyl. C ₁₇ H ₂₄ O ₉ 372.37 USP Eleutheroside E RS ?-D-Glucopyran oside, (tet rahydr o-1 <i>H</i> ,3 <i>H</i> -furo(3,4-c)furan -1,4-diyl)bis(2,6- dimethoxy-4,1-p henylene)bis-. C ₃₄ H ₄₆ O ₁₈ 742.70 to: USP Eleutheroside B RS USP Eleutheroside E RS
ESTROPIPATE CHEMICAL INFORMATION	USPNF 2021 ISSUE 1	Online	19-Nov-2021	1-Dec-2021	NA	NA	Change 436.56 to: 436.57
ISOSORBIDE DINITRATE EX <711>	USPNF 2021 ISSUE 1	Online	19-Nov-2021	1-Dec-2021	NA	NA	In Test 2: Change

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TENDED- RELEASE TABLETS								Determine the amount of isosorbide dinitrate ($C_6H_9NO_6$) dissolved by employing the following method. to: Determine the amount of isosorbide dinitrate ($C_6H_8N_2O_8$) dissolved by employing the following method.
APREPITANT CAPSULES	PERFORMANC E TESTS/ <i>Dissolution</i> <711>	USPNF 2021 ISSUE 1	Online	19-Nov-2021	1-Dec-2021	NA	NA	In <i>Test 1/Apparatus 2</i> : Change [Note—A suitable sinker is available from VanKel, www.chem.agilent.com , catalog number 12-3050. Proper placement of

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ISOPROTERE NOL HYDROC HLORIDE INJECTION	<i>Color and clarity</i>	USPNF 2021 <i>ISSUE 1</i>	Online	19-Nov-2021		1-Dec-2021	NA	NA	the Capsules is in the sinkers with the cap facing the fixed prong end.] to: [Note—A suitable sinker is available from www.agilent.com, catalog number 12-3050. Proper placement of the Capsules is in the sinkers with the cap facing the fixed prong end.] Change Using the Injection as the <i>Test solution</i>, proceed as directed for <i>Color and clarity</i> under <i>Isoproterenol Inhalation Solution</i>. to:

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							<i>Standard solution</i> —Transfer 2.0 mL of 0.100 N iodine VS to a 500-mL volumetric flask, dilute with water to volume, and mix. <i>Procedure</i>—Visually examine a portion of the Injection (<i>Test solution</i>) in a suitable clear glass test tube against a white background: it is not pinkish and it contains no precipitate. If any yellow color is observed in the <i>Test solution</i>, concomitantly determine the absorbances of the <i>Test</i>

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ELEUTHERO ROOT AND RHIZOME POWDER	ADDITIONAL REQUIREMENT S/USP Reference Standards <11>	USPNF 2021 ISSUE 1	Online	19-Nov-2021		1-Dec-2021	NA	NA	<i>solution</i> and the <i>Standard solution</i> in 1-cm cells with a suitable spectrophotometer set at 460 nm: the absorbance of the <i>Test solution</i> does not exceed that of the <i>Standard solution</i>. Change USP Eleutheroside B RS ?-D-Glucopyranoside, 4-(3-hydroxy-1-propenyl)-2,6-dimethoxyphenyl. $C_{17}H_{24}O_9$ 372.37 USP Eleutheroside E RS ?-D-Glucopyranoside, (tetrahydr

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							o-1<i>H</i>,3<i>H</i>-furo(3,4-c)furan-1,4-diyl)bis(2,6-dimethoxy-4,1-p henylene)bis-. C₃₄H₄₆O₁₈ 742.70 to: USP Eleutheroside B RS USP Eleutheroside E RS
DACARBAZINE IM PUR ITIES/ <i>Organic Impurities</i>	USPNF 2021 ISSUE 1	Online	19-Nov-2021	1-Dec-2021	NA	NA	In <i>System suitability/Suitability requirements/Sig nal-to-noise ratio</i>: Change NTL 10 to: NLT 10
CARBOMER IN IM TERPOLYMER PURITIES/ <i>Limit of Benzene</i>	USPNF 2021 ISSUE 2	Online	19-Nov-2021	1-Dec-2021	NA	NA	In <i>Analysis</i>: Change W_U = weight of Carbomer Interpolymer in the <i>Sample solution</i> (mg) to: W

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CEFTIOFUR H YDROCHLORIDE	IM PURITIES/Low Molecular Weight Impurities	USPNF 2021 ISSUE 1	Online	19-Nov-2021		1-Dec-2021	NA	NA	U = weight of Carbomer Interpolymer in the Sample solution (?g) In Analysis: Change Result = $\{r_U/[r_S + (?r_U/F)]\} \times (1/F) \times 100$ to: Result = $\{r_U/[r_S + ?(r_U/F)]\} \times (1/F) \times 100$
DOXAZOSIN MESYLATE	ADDITIONAL REQUIREMENT S/USP Reference Standards <11>	USPNF 2021 ISSUE 1	Online	29-Oct-2021		1-Nov-2021	NA	NA	In USP Doxazosin Related Compound D RS: Change 1,4-Benzodioxane-2-carboxylic acid. $C_9H_8O_5$ 196.16 to: 1,4-Benzodioxane-2-carboxylic acid. $C_9H_8O_4$ 180.16 Change Result = $[(C/V)/W] \times 100$
ECHOTHIOPHATE IODIDE FOR	SPECIFIC TESTS/ Water/Analysis	USPNF2021 Issue 1	Online	29-Oct-2021		1-Nov-2021	NA	NA	Change Result = $[(C/V)/W] \times 100$

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OPHTHALMIC SOLUTION							to: Result = [(C × V)/W] × 100
TIAGABINE HYDROCHLORIDE ADDITIONAL REQUIREMENTS	USPNF 2021 ISSUE 1 S/USP Reference Standards <11>	Online	29-Oct-2021	1-Nov-2021	NA	NA	In USP Tiagabine Hydrochloride RS: Delete [NOTE—Available in monohydrate form.]
DIMENHYDRINATE ORAL SOLUTION	Content of 8-chlorotheophylline USPNF 2021 ISSUE 1	Online	29-Oct-2021	1-Nov-2021	NA	NA	Change Ammonium bicarbonate solution, Diluent, Solution A, Solution B, Mobile phase, Internal standard solution, and Chromatographic system —Proceed as directed in the Assay under Dimenhydrinate Tablets. to:

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							<i>Ammonium bicarbonate solution</i> —Dissolve 4 g of ammonium bicarbonate in 250 mL of water. <i>Diluent</i> —Dissolve 4 g of ammonium bicarbonate in 200 mL of water. Add 50 mL of methanol, and mix. <i>Solution</i> <i>A</i>—Dissolve 0.8 g of ammonium bicarbonate in 800 mL of water. Add 200 mL of methanol, filter, and degas. <i>Solution</i> <i>B</i>—Dissolve 0.8 g of ammonium bicarbonate in 150 mL of water. Add 850

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							mL of methanol, filter, and degas. <i>Mobile phase</i> —Use variable mixtures of <i>Solution A</i> and <i>Solution B</i> as directed for <i>Chromatographic system</i> . Make adjustments if necessary (see <i>System Suitability</i> under <i>Chromatography</i> ?621?). <i>Internal standard solution</i> —Prepare a solution in methanol containing 2.0 mg of 2-hydroxy benzyl alcohol per mL. AND Change <i>Standard</i>

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							<i>solution</i> —Prepare as directed for <i>Standard preparation</i> in the Assay under <i>Dimenhydrinate Tablets</i>. to: <i>Standard solution</i> —Accurately weigh about 50 mg of USP Dimenhydrinate RS, add about 5 mL of <i>Ammonium bicarbonate solution</i> and 20.0 mL of <i>Internal standard solution</i>, and mix. To 1 mL of this solution add about 9 mL of <i>Diluent</i>, and mix. AND Add

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							<i>Chromatographic system (see Chromatography ?621?)—The liquid chromatograph is equipped with a 229-nm detector and a 4.6-mm × 25-cm column that contains packing L7. The flow rate is about 1.5 mL per minute. The chromatograph is programmed as follows.</i> Time (minutes), 0, 0–7.0, 7.0–7.1, 7.1–15, 15–15.1, 15.1–22.0 Solution A (%), 100, 100, 100?0, 0, 0?100, 100 Solution B (%), 0, 0, 0?100,

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							100, 100?0, 0 Elution, equilibration, isocratic, linear gradient, isocratic, linear gradient, isocratic
							Chromatograph the <i>Standard solution</i> , and record the peak areas as directed for <i>Procedure</i> : the relative retention times are about 0.3 for 8-chlorotheo phylline, 0.5 for the internal standard, and 1.0 for diphenhy dramine; the resolution, <i>R</i> , between 8-chlor otheophylline and the internal standard is not less than 4.5;

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DOXAZOSIN MESYLATE	IMPURITIES/ <i>Organic Impurities</i>	USPNF 2021 ISSUE 1	Online	29-Oct-2021		1-Nov-2021	NA	NA	and the relative standard deviation for replicate injections is not more than 2.0%. In <i>Table 2</i> : Change Doxazosin related compound D ^g 0.83, 196.16, 196.16, 0.25 to: Doxazosin related compound D ^g 0.83, 180.16, 180.16, 0.25
ARTICLES OF BOTANICAL ORIGIN	PESTICIDE RESIDUE ANALYSIS/ <i>Limits</i>	USPNF2021 Issue 1	Online	29-Oct-2021		1-Nov-2021	NA	NA	In <i>Qualitative and Quantitative Analysis of Pesticide Residues</i> : Change [Note—Current version Document No. SANTE/11813/2

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ALFUZOSIN H IM YDROCHLORI PUR DE EXTENDEDITIES/ <i>Organic</i> RELEASE <i>Impurities</i> TABLETS	USPNF 2021 ISSUE 1	Online	29-Oct-2021	1-Nov-2021	NA	NA	017, https://ec.europa.eu/food/sites/food/files/plant/docs/pesticides_mrl_guidelines_wrkdoc_2017-1813.pdf to: [Note—Current version Document No. SANTE/12682/2019, https://ec.europa.eu/food/system/files/2020-01/pesticides_mrl_guidelines_wrkdoc_2019-12682.pdf] In Analysis: Change r_S = peak response of alfuzosin from the Sample solution to: r_S = peak response of

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DIMENHYDRIN Assay ATE ORAL SOLUTION	<i>USPNF 2021 Issue 1</i>	Online	29-Oct-2021	1-Nov-2021	NA	NA	alfuzosin from the <i>Standard solution</i> Change <i>Ammonium bicarbonate solution, Diluent, Solution A, Solution B, Mobile phase, Internal standard solution, Standard preparation, and Chromatographic system</i> —Proceed as directed in the Assay under <i>Dimenhydrinate Tablets</i>. to: <i>Ammonium bicarbonate solution</i> —Dissolve 4 g of ammonium

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							bicarbonate in 250 mL of water. <i>Diluent</i> —Dissolve 4 g of ammonium bicarbonate in 200 mL of water. Add 50 mL of methanol, and mix. <i>Solution</i> <i>A</i> —Dissolve 0.8 g of ammonium bicarbonate in 800 mL of water. Add 200 mL of methanol, filter, and degas. <i>Solution</i> <i>B</i> —Dissolve 0.8 g of ammonium bicarbonate in 150 mL of water. Add 850 mL of methanol, filter, and degas. <i>Mobile phase</i> —Use

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							variable mixtures of <i>Solution A</i> and <i>Solution B</i> as directed for <i>Chromatographic</i> system. Make adjustments if necessary (see <i>System Suitability</i> under <i>Chromatography</i> ?621?). <i>Internal standard solution</i> —Prepare a solution in methanol containing 2.0 mg of 2-hydroxy benzyl alcohol per mL. <i>Standard preparation</i> —Accurately weigh about 50 mg of USP Dimenhydrinate RS, add about 5

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							mL of <i>Ammonium bicarbonate solution</i> and 20.0 mL of <i>Internal standard solution</i>, and mix. To 1 mL of this solution add about 9 mL of <i>Diluent</i>, and mix. AND Add <i>Chromatographic system</i> (see <i>Chromatography</i> ?621?)—The liquid chromatograph is equipped with a 229-nm detector and a 4.6-mm × 25-cm column that contains packing L7. The flow rate is about 1.5 mL per minute. The

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							chromatograph is programmed as follows. Time (minutes), 0, 0–7.0, 7.0–7.1, 7.1–15, 15–15.1, 15.1–22.0 Solution A (%), 100, 100, 100?0, 0, 0?100, 100 Solution B (%), 0, 0, 0?100, 100, 100?0, 0 Elution, equilibration, isocratic, linear gradient, isocratic, linear gradient, isocratic Chromatograph the <i>Standard preparation</i>, and record the peak areas as directed for <i>Procedure</i>: the

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							relative retention times are about 0.3 for 8-chlorotheophylline, 0.5 for the internal standard, and 1.0 for diphenhydramine; the resolution, <i>R</i>, between 8-chlorotheophylline and the internal standard is not less than 4.5; and the relative standard deviation for replicate injections is not more than 2.0%. AND Change <i>Procedure</i>—Proceed as directed for <i>Procedure</i> in the Assay under <i>Dimenhydrinate Tablets</i>.

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PAROXETINE CHEMICAL HYDROCHLOR INFORMATION IDE	<i>Interim Revision Announcement (Official May 01, 2021)</i>	Online	29-Oct-2021	1-Nov-2021	NA	NA	to: <i>Procedure</i> —Separately inject equal volumes (about 10 µL) of the <i>Standard preparation</i> and the <i>Assay preparation</i> into the chromatograph, record the chromatograms, and measure the areas for the major peaks. Change (?)-(3<i>S</i>,4<i>R</i>)-4-(<i>p</i>-Fluorophenyl)-3-[(3,4-methylenedioxy)phenoxy]methyl]piperidine hydrochloride; to: (?)-(3<i>S</i>,4<i>R</i>)-4-(<i>p</i>

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SACCHARIN	SPECIFIC TESTS/ <i>Readily Carbonizable Substances Test <271></i>	USPNF 2021 ISSUE 1	Online	24-Sep-2021	1-Oct-2021	NA	NA	-Fluorophenyl)-3-[[[(3,4-methylenedioxy)phenoxy)methyl]piperidine hydrochloride; In <i>Matching fluid A</i> : Change Cobaltous chloride CS, ferric chloride TS, cupric sulfate CS, and water (0.1:0.4:0.1:4.4) to: Cobaltous chloride CS, ferric chloride CS, cupric sulfate CS, and water (0.1:0.4:0.1:4.4)
GLYBURIDE TABLETS	ADDITIONAL REQUIREMENT <i>SI/USP Reference Standards <11></i>	USPNF 2021 ISSUE 1	Online	24-Sep-2021	1-Oct-2021	NA	NA	In USP Glyburide Related Compound A RS: Change 4-[2-(5-Chloro-2-methoxybenzamide)ethyl]ben

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ACETAMINOPHEN TABLETS PURITIES/Organic Impurities	USPNF 2021 ISSUE 1	Online	24-Sep-2021	1-Oct-2021	NA	NA	zenesulfonamide. to: 4-[2-(5-Chloro-2-methoxybenzamide)ethyl]benzenesulfonamide. In <i>Chromatographic system/ Detector:</i> Change UV 272 nm. For <i>Identification B</i> , use a diode array detector in the range of 200–400 nm. to: UV 272 nm In <i>Table 1:</i> Change ?1430.6? ^a to: ?1430.6? AND Change ?1430.7? ^a to:
ANALYTICAL METHODS BASED ON SCATTERING PHENOMENA—SCATTERING GENERAL	1. OVERVIEW: GENERAL BASED ON SCATTERING PHENOMENA USPNF 2021 ISSUE 1	Online	24-Sep-2021	1-Oct-2021	NA	NA	

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LIGHT OBSCURATION METHOD FOR THE DETERMINATION OF SUBVISIBLE PARTICULATE MATTER	USPNF 2021 ISSUE 1	Online	24-Sep-2021	1-Oct-2021	NA	NA	?1430.7? AND Delete footnote a In <i>Overview</i> : Change The test is also sensitive to size accuracy at 15 ?m, since small errors in size will translate into large errors in count at the midpoint of the PS distribution. to: The test is also sensitive to size accuracy at 15 ?m, since small errors in size will translate into large errors in count at the midpoint of the particle size distribution. In <i>Analysis</i> : Change To 2.0 mL of
HYDROXYETHYL CELLULOSE IM PUR ITIES/	USPNF 2021 ISSUE 1	Online	24-Sep-2021	1-Oct-2021	NA	NA	

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									the <i>Sample solution</i>, add 5.0 mL of a 4-g/L solution of methylbenzothiazolone hydrazone hydrochloride to an 80% (v/v) solution of glacial acetic acid in water. to: To 2.0 mL of the <i>Sample solution</i>, add 5.0 mL of a 4-g/L solution of methylbenzothiazolone hydrazone hydrochloride in an 80% (v/v) solution of glacial acetic acid in water.
OFLOXACIN	CHEMICAL INFORMATION	USP NF 2021 ISSUE 1	Online	24-Sep-2021		1-Oct-2021	NA	NA	Change (±)-9-Fluoro-2,3-dihydro-3-methyl-10-(4-methyl-1-piperazinyl)-7-o

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ANALYTICAL METHODS BASED ON SCATTERING PHENOMENA-GENERAL	2. INTRODUCT IES	USPNF 2021 ISSUE 1	Online	24-Sep-2021		1-Oct-2021	NA	NA	xo-7H -p yrido [1,2,3-de]-1,4-benzoxazi ne-6-carboxylic acid to: (±)-9-Fluoro-2, 3-dihydro-3-met hyl-10-(4-methyl -1-piperaziny)-7 -oxo-7H -p yrido [1,2,3-de]-1,4-benzoxazi ne-6-carboxylic acid In paragraph five: Change <i>Determination of Zeta Potential by Electrophoretic Light Scattering</i> ?432?¹ to: <i>Determination of Zeta Potential by</i>

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PYRANTEL PAMOATE	ADDITIONAL REQUIREMENTS/USP Reference Standards <11>	USPNF 2021 ISSUE 2	Online	24-Sep-2021		1-Oct-2021	NA	NA	<i>Electrophoretic Light Scattering</i> ?432? AND In footnote 1: Change This chapter will appear in a future <i>Pharmacopeial Forum (PF)</i> issue. to: This chapter appeared in issue 46(3) of the <i>Pharmacopeial Forum (PF)</i>. In USP Pyrantel Related Compound A RS: Change (Z)-1-Methyl-2-(2-(thiophen-2-yl)vinyl)-1,4,5,6-tetrahydropyrimidine. C₃₄H₃₀N₂O₆S 594.69

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RIFABUTIN	IMPURITIES/Organic Impurities, Procedure 2	USPNF 2021 ISSUE 1	Online	27-Aug-2021		1-Sep-2021	NA	NA	to: (Z)-1-Methyl-2-[2-(2-thiophenyl)vinyl]-1,4,5,6-tetrahydropyrimidine 4,4'-methylene bis[3-hydroxy-2-naphthoate] (1:1). C₁₁H₁₄N₂S · C₂₃H₁₆O₆ 594.68 In <i>Sample solution</i>: Change 1 mg/mL of USP Rifabutin RS prepared as follows. to: 1 mg/mL of Rifabutin prepared as follows.
SPECTROSCOPIC IDENTIFICATION TESTS	EQUIVALENT/ALTERNATIVE TESTS	USPNF 2021 ISSUE 1	Online	27-Aug-2021		1-Sep-2021	NA	NA	Change (see <i>Mass Spectrometry</i> ?736?, <i>Nuclear Magnetic Resonance</i>

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							<i>Spectroscopy</i> ?761?, ?854?, ?857?, ?941?, <i>Near-Infrared S</i> <i>pectroscopy—Th</i> <i>eory and</i> <i>Practice</i> ?1856?, and ?858?. to: (see <i>Mass</i> <i>Spectrometry</i> ?736?, <i>Nuclear</i> <i>Magnetic</i> <i>Resonance</i> <i>Spectroscopy</i> ?761?, ?854?, ?856?, ?857?, ?858?, and ?941?).

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