

**ERRATA TO FCC 6TH EDITION 1ST Supplement
Version 1 June 2009**

Date Posted	FCC 6, 1S Page	Title	Section	Description
6/2009	1220	Annotated List	<i>Manganese Sulfate</i>	Change “ <i>Impurities</i> (section <i>Organic impurities</i>)” to “ <i>Impurities</i> (section <i>Inorganic impurities</i>)”
6/2009	1273	Food Starch, Modified	OTHER REQUIREMENTS/ Monofunctional and/or Polyfunctional Esterification (Starch Esters)/Treatment to Produce Starch Acetate	Change “Acetic anhydride of vinyl acetate” to “Acetic anhydride or vinyl acetate”
6/2009	1325	Pullulan	CAS	Change “[9054-02-07]” to “[9057-02-07]”
6/2009	1355	Appendix II: Physical Tests and Determinations	A. CHROMATOGRAPHY/ HIGH-PRESSURE LIQUID CHROMAGOTRAPHY/ Interpretation of Chromatograms	Page 1364, line 21: Change “X is the mean of the set of N measurements” to “ \bar{X} is the mean of the set of N measurements”
6/2009	1384	Appendix III: Chemical Tests and Determinations	C. OTHERS/COLORS/ Total Color	Page 1412, lines 18, 19, 20 and 25: Change “ $m = [N\sum C_i A_i - \sum C_i \sum A_i] / [N\sum A_i^2 - (\sum A_i)^2]$, $b = [A]_i - m[C]_i$, in which C and A are the calculated averages of the concentrations and peak areas, respectively, used to construct the standard curve for one intermediate or side reaction product. Calculate the correlation coefficient, r, from the following equation: $r = [\sum (C_i - C)(A_i - A)] / [\sum (C_i - C)^2 \times \sum (A_i - A)^2]$.” to “ $m = [N\sum C_i A_i - \sum C_i \sum A_i] / [N\sum A_i^2 - (\sum A_i)^2]$, $b = [\bar{A}]_i - m[\bar{C}]_i$ in which \bar{C} and \bar{A} are the calculated averages of the

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				<p>concentrations and peak areas, respectively, used to construct the standard curve for one intermediate or side reaction product. Calculate the correlation coefficient, r, from the following equation:</p> $r = \frac{[\sum(C_i - \bar{C})(A_i - \bar{A})]}{[\sum(C_i - \bar{C})^2 \times \sum(A_i - \bar{A})^2]}.$
6/2009	1384	Appendix IIIC. OTHERS	APPENDIX II C. OTHERS / Benzene/Calibration	In the Calculation for the response factor for benzene, change " B_v " to " W_b "