



# SPECIFICATIONS



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ANIMAL NUTRITION

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Salisbury, MD 21804

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| Standard Analysis | % DM |
|-------------------|------|
| Moisture          | 6.3  |
| Dry Matter        | 93.7 |
| Nitrogen          | 14.8 |
| Crude Protein     | 92.3 |
| Soluble Protein   | 41.8 |
| Ammonia           | 0.0  |
| NDIP              | 8.2  |
| ADIP              | 8.2  |
| Crude Fiber       | 0.0  |
| ADF               | 0.0  |
| NDF               | 0.0  |
| Lignin            | 0.0  |
| Total VFA         | 0.0  |
| Lactic            | 0.0  |
| Acetic            | 0.0  |
| Sugar             | 0.0  |
| Starch            | 0.0  |
| Soluble Fiber     | 0.0  |
| NFC               | 0.0  |

| Protein Bioavailability | % DM |
|-------------------------|------|
| RUP % CP                | 93.3 |
| RUPd % RUP              | 90.5 |
| DRUP % DM               | 77.9 |
| DRUP-Lys % DM           | 6.6  |
| DRUP-Met % DM           | 4.3  |
| DRUP-EAA % DM           | 45.1 |

| Macro Mineral | % DM |
|---------------|------|
| Ash           | 5.4  |
| Calcium       | 0.6  |
| Phosphorus    | 0.4  |
| Magnesium     | 0.0  |
| Potassium     | 0.2  |
| Sulfur        | 0.7  |
| Sodium        | 0.4  |
| Chloride      | 0.5  |

| Trace Mineral | ppm   |
|---------------|-------|
| Iron          | 2,323 |
| Manganese     | 4.7   |
| Zinc          | 35    |
| Copper        | 6.5   |

| Fatty Acids    | % DM |
|----------------|------|
| Fat            | 2.3  |
| TFA            | 2.2  |
| Glycerol       | 0.1  |
| Pigment        | 0.1  |
| C12:0          | 0.0  |
| C14:0          | 0.0  |
| C16:0          | 0.6  |
| C16:1          | 0.1  |
| C18:0          | 0.7  |
| C18:1T         | 0.0  |
| C18:1C         | 0.5  |
| C18:2          | 0.2  |
| C18:3          | 0.0  |
| Other          | 0.1  |
| Fat Type       | 1.9  |
| Lipolysis Rate | 55.7 |

| Pool   | % DM  | Rate | Int dig |
|--------|-------|------|---------|
| CHO A1 | -     | -    | -       |
| CHO A4 | -     | -    | -       |
| CHO B1 | -     | -    | -       |
| CHO B2 | -     | -    | -       |
| CHO B3 | -     | -    | -       |
| CHO C  | -     | -    | -       |
| PRO A1 | -     | -    | -       |
| PRO A2 | 41.82 | 1.00 | 100.0   |
| PRO B1 | 42.34 | 0.55 | 100.0   |
| PRO B2 | -     | -    | -       |
| PRO C  | 8.19  | -    | -       |

| Amino Acids   | Product |         | Model Specifications <sup>1</sup> |         |
|---------------|---------|---------|-----------------------------------|---------|
|               | % DM    | % of CP | % DM                              | % of CP |
| Arginine      | 3.7     | 4.0     | 3.6                               | 3.9     |
| Histidine     | 5.2     | 5.7     | 5.5                               | 6.0     |
| Isoleucine    | 1.0     | 1.0     | 0.8                               | 0.9     |
| Leucine       | 11.6    | 12.6    | 11.9                              | 12.9    |
| Lysine        | 7.6     | 8.2     | 7.8                               | 8.4     |
| Methionine    | 4.9     | 5.3     | 5.1                               | 5.5     |
| Phenylalanine | 6.2     | 6.7     | 6.3                               | 6.9     |
| Threonine     | 3.7     | 4.0     | 3.8                               | 4.1     |
| Tryptophan    | 1.1     | 1.2     | 1.1                               | 1.2     |
| Valine        | 7.6     | 8.2     | 7.7                               | 8.3     |
| EAA           | 52.7    | 57.0    | 53.4                              | 57.9    |

<sup>1</sup>When a feed component of a dairy ration consists of a blend of feed ingredients differing in rumen undegradable protein and/or post-ruminal protein digestibility, the amino acids of the resulting mixture differ in their ruminal degradability and post ruminal digestibility. Most ration formulation software allow only a single, common set of parameters to characterize ruminal degradability and post ruminal digestibility of all essential amino acids; one cannot enter parameters specific to each amino acid. Therefore, entering the amino acid profile of the whole protein of a feed mixture results in erroneous values of its metabolizable amino acid content. To circumvent this problem, users should use the 'model specifications' amino acid values in software such as CNCPS/NASEM so that the correct supply of each metabolizable amino acid is calculated correctly.