

## Molecular Weight/Isoelectric Point

### 1. Purpose

This document describes the process by which protein parameters are calculated, involving isoelectric point (pI) and molecular weight (MW). Values are calculated by a BioPerl related script, in a precompute process using GenBank files, and are presented to the user in VIPR.

### 2. Method Description

The method calculates physical parameters of proteins given only their sequence. All proteomics studies that separate proteins on the basis of their charge and size generate quantitative values. By calculating these values in a precompute, users can search for molecules within a given range and size, to help determine what they have been studying and purifying.

### 3. Input Data Preparation

Input to the method is a protein fasta (.faa) amino acid text file, such as the protein sequence:

```
>PB1_F2
MEQEQDTPWTQSTGHINIQKRGNGQQTQRLLEHLNSTRLTGHCLRTMSQVDMHKQTVSWKQW
LSLKSPTQESLKTRVLKRWKLSNKQEWTN
```

### 4. Output Data, Processing and Display

Intermediate output results are of the delimited format:

gi	iep	mw_lower	mw_upper	diff
110669658	8.71	57676.4	57676.4	t
110669659	5.11	41674.3	41674.3	t
110669660	6.68	30305.3	30951.1	f

The lower and upper MW difference occurs ('f') if a sequencing abnormality reveals an amino acid residue whose identity cannot be determined precisely ('X'). For example, a string of four X's can result in the calculation of MW that differ in lower and upper range by 645 g/mol (i.e. all glycine vs. all tryptophan).

Results such as for the NA fragment above are presented back to the user as shown:

Isoelectric pt/Molecular Weight ([SOP](#))

Isoelectric pt	Molecular Weight	Evidence Code
6.9	255972.8	RCA

The precision of the method is pI is +/- 0.01 charge unit. MW precision is to the closest 0.01 g/mol.

### 5. Reference

1. Stajich, J. et al. (2002) The BioPerl Toolkit: Perl Modules for the Life Sciences. *Genome Res.* 12(10): 1611-1618. URL: <http://doc.BioPerl.org/releases/BioPerl-1.0.1/Bio/SeqIO.html>