

**NAME**

ElementalAnalysis.pl - Perform elemental analysis using specified formulas

**SYNOPSIS**

ElementalAnalysis.pl Formula(s)...

```
ElementalAnalysis.pl [-h, --help] [-m, --mode All | "ElementalAnalysis, [MolecularWeight, ExactMass]"] [
--outdelim comma | tab | semicolon] [--output STDOUT | File] [--outputstyle FormulaBlock | FormulaRows] [
-o, --overwrite] [--precision number] [-q, --quote yes | no] [-r, --root rootname] [-v --valueLabels [Name,
Label, [Name, Label,...]] [-w, --workingdir dirname] Formula(s)...
```

**DESCRIPTION**

Perform elemental analysis using molecular formula(s) specified on the command line.

In addition to straightforward molecular formulas - H<sub>2</sub>O, HCl, C<sub>3</sub>H<sub>7</sub>O<sub>2</sub>N - other supported variations are: Ca<sub>3</sub>(PO<sub>4</sub>)<sub>2</sub>, [PCl<sub>4</sub>]<sup>+</sup>, [Fe(CN)<sub>6</sub>]<sup>4-</sup>, C<sub>3</sub>H<sub>4</sub>N<sub>2</sub>O<sub>6</sub>+<sub>2</sub>, Na<sub>2</sub>CO<sub>3</sub>·10H<sub>2</sub>O, 8H<sub>2</sub>S·46H<sub>2</sub>O, and so on. Charges are simply ignored. Isotope symbols in formulas specification, including D and T, are not supported.

**PARAMETERS**

Formulas *Formula1* [*Formula2*...]

*Formulas* is a space delimited list of molecular formulas to use for elemental analysis.

Input value format is: *Formula1* [*Formula2* *Formula3*...]. Default: *H<sub>2</sub>O*. Examples:

```
HCl
HCl, C3H7O2N
H2O2 Ca3(PO4)2 [PCl4]+
```

**OPTIONS**

-h, --help

Print this help message.

--fast

In this mode, the specified formulas are considered valid and initial formula validation check is skipped.

-m, --mode *All* | "*ElementalAnalysis*,*[MolecularWeight,ExactMass]*"

Specify what values to calculate using molecular formulas specified on command line: calculate all supported values or specify a comma delimited list of values. Possible values: *All* | "*ElementalAnalysis*,*[MolecularWeight, ExactMass]*". Default: *All*.

--outdelim *comma* | *tab* | *semicolon*

Output text file delimiter. Possible values: *comma*, *tab*, or *semicolon* Default value: *comma*.

--output *STDOUT* | *File*

List information at *STDOUT* or write it to a file. Possible values: *STDOUT* or *File*. Default: *STDOUT*. -r, --root option is used to generate output file name.

--outputstyle *FormulaBlock* | *FormulaRows*

Specify how to list calculated values: add a new line for each property and present it as a block for each formula; or include all properties in one line and show it as a single line.

Possible values: *FormulaBlock* | *FormulaRows*. Default: *FormulaBlock*

An example for *FormulaBlock* output style:

```
Formula: H2O
ElementalAnalysis: H: H: 11.1898%; O: 88.8102%
MolecularWeight: 18.0153
ExactMass: 18.0106
... ..
... ..
... ..
```

```

Formula: H2O2
ElementalAnalysis: H: 5.9265%; O: 94.0735%
MolecularWeight: 34.0147
ExactMass: 34.0055
... ..
... ..
... ..

```

An example for *FormulaRows* output style:

```

Formula,ElementalAnalysis,MolecularWeight,ExactMass
H2O,H: 11.1898%; O: 88.8102%,18.0153,18.0106
H2O2,H: 5.9265%; O: 94.0735%,34.0147,34.0055

```

**-o, --overwrite**

Overwrite existing files.

**--precision *number***

Precision for listing numerical values. Default: up to 4 decimal places. Valid values: positive integers.

**-r, --root *rootname***

New text file name is generated using the root: <Root>.<Ext>. File name is only used during *File* value of **-o, --output option**.

Default file name: FormulsElementalAnalysis.<Ext>. The csv, and tsv <Ext> values are used for comma/semicolon, and tab delimited text files respectively.

**-v --valueLabels *Name,Label,[Name,Label,...]***

Specify labels to use for calculated values. In general, it's a comma delimited list of value name and column label pairs. Supported value names: *ElementalAnalysis, MolecularWeight, and ExactMass*. Default labels: *ElementalAnalysis, MolecularWeight, and ExactMass*.

**-w, --workingdir *dirname***

Location of working directory. Default: current directory.

## EXAMPLES

To perform elemental analysis, calculate molecular weight and exact mass for H2O, type:

```
% ElementalAnalysis.pl
```

To perform elemental analysis, calculate molecular weight and exact mass for Ca3(PO4)2 and [PCl4]+, type:

```
% ElementalAnalysis.pl "Ca3(PO4)2" "[PCl4]+"
```

To perform elemental analysis, use label analysis for calculated data, and generate a new CSV file ElementalAnalysis.csv for H2O and H2O2, type:

```
% ElementalAnalysis.pl --m ElementalAnalysis --output File
--valueLabels "ElementalAnalysis,Analysis" -o -r ElementalAnalysis.csv
H2O H2O2
```

To calculate molecular weight and exact mass with four decimal precision and generate a new CSV file WeightAndMass.csv with data rows for H2O and H2O2, type:

```
% ElementalAnalysis.pl --m "MolecularWeight,ExactMass" --output File
--outputstyle FormulaRows -o -r WeightAndMass.csv
H2O H2O2
```

## AUTHOR

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SEE ALSO

[ElementalAnalysisSDFFiles.pl](#), [ElementalAnalysisTextFiles.pl](#)

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