

Edit parameters for Align with MafftWS



Current parameter set name : My Defaults Revised

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Details

description

Options

Shared 6mers distance calculation

Output sequences order

--inputorder

Sequence type

--auto

Pairwise alignment computation...

--localpair

--nuc - Assume the sequences are nucleotide. --amino - Assume the sequences are amino acid.

[PLink](#)

Parameters

Max iteration number



0

Partsize

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