

Βούρδουλα Καλλιόπη-Αικατερίνη  
8289

### Πρωτεΐνη: 3LYM(P00698)

- Fasta sequence:

```
>sp|P00698|LYSC_CHICK Lysozyme C OS=Gallus gallus  
GN=LYZ PE=1 SV=1  
MRSLLILVLCFLPLAALGKVFGRCELAAMKRHGLDNYRGYSLGNWVCAAKFE  
SNFNTQA  
TNRNTDGSTDYGILQINSRWWCNDGRTPGSRNLCNIPCSALLSSDITASVNCA  
KKIVSDG  
NGMNAWVAWRNRCKGTDVQAWIRGCRL
```

- ProtParam

#### User-provided sequence:

```
      10           20           30           40  
50 MRSLILVLC FLPLAALGKV FGRCELAAM KRHGLDNYRG  
YSLGNWVCAA KFESNFNTQA  
  
      70           80           90          100  
110 TNRNTDGSTD YGILQINSRW WCNDGRTPGS RNLNIPCSA  
LLSSDITASV NCAKKIVSDG  
  
      130          140  
NGMNAWVAWR NRCKGTDVQA WIRGCRL
```

[References](#) and [documentation](#) are available.

**Number of amino acids:** 147

**Molecular weight:** 16238.6

**Theoretical pI:** 9.36

**Amino acid composition:**

Ala (A)	14	9.5%
Arg (R)	12	8.2%
Asn (N)	14	9.5%
Asp (D)	7	4.8%
Cys (C)	9	6.1%
Gln (Q)	3	2.0%
Glu (E)	2	1.4%
Gly (G)	13	8.8%
His (H)	1	0.7%
Ile (I)	7	4.8%
Leu (L)	15	10.2%
Lys (K)	6	4.1%
Met (M)	3	2.0%
Phe (F)	4	2.7%
Pro (P)	3	2.0%
Ser (S)	11	7.5%
Thr (T)	7	4.8%
Trp (W)	6	4.1%
Tyr (Y)	3	2.0%
Val (V)	7	4.8%
Pyl (O)	0	0.0%
Sec (U)	0	0.0%

(B)	0	0.0%
(Z)	0	0.0%
(X)	0	0.0%

**Total number of negatively charged residues (Asp + Glu): 9**

**Total number of positively charged residues (Arg + Lys): 18**

**Atomic composition:**

Carbon	C	705
Hydrogen	H	1116
Nitrogen	N	214
Oxygen	O	204
Sulfur	S	12

**Formula:** C<sub>705</sub>H<sub>1116</sub>N<sub>214</sub>O<sub>204</sub>S<sub>12</sub>

**Total number of atoms:** 2251

**Extinction coefficients:**

Extinction coefficients are in units of  $M^{-1} \text{ cm}^{-1}$ , at 280 nm measured in water.

Ext. coefficient      37970  
Abs 0.1% (=1 g/l)    2.338, assuming all pairs of Cys residues form cystines

Ext. coefficient      37470  
Abs 0.1% (=1 g/l)    2.307, assuming all Cys residues are reduced

**Estimated half-life:**

The N-terminal of the sequence considered is M (Met).

The estimated half-life is: 30 hours (mammalian reticulocytes, in vitro).

>20 hours (yeast, in vivo).

>10 hours (Escherichia coli, in vivo).

**Instability index:**

The instability index (II) is computed to be 19.86  
This classifies the protein as stable.

**Aliphatic index:** 81.70

**Grand average of hydropathicity (GRAVY):** -0.150

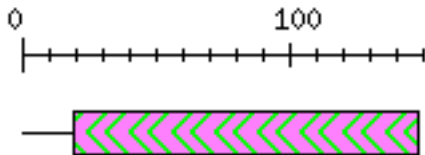
- ProDom

<b>database:</b>	multiple alignments
<b>Program:</b>	ncbi-blastp
<b>Matrix:</b>	BLOSUM62

**Expect:** 0.01  
**Filter:** seg

## Graphical results and forms to other applications

The following is the graphical representation of the HSP found by BLAST. Please note that HSPs are sorted from highest to lowest scores, so that lower scoring HSPs may be hidden.



## Align subsequence with ProDom domains, using Multalin

Domain ID	BEGIN	END
PD00057 7	19	147
PDC5E2N 9	45	145
PDC1J6J 2	42	134

## Domain 3D modelling using Swiss-Model

Domain ID	BEGIN	END
PD00057 7	19	147

## Domain 3D modelling using Geno3D

Domain ID	BEGIN	END
PD000577	19	147

## HSP Results

Warning: Original output has been filtered to yield non-redundant similarities

blastp 2.2.26 [Sep-21-2011]

Reference: Altschul, Stephen F., Thomas L. Madden, Alejandro A. Schaffer, Jinghui Zhang, Zheng Zhang, Webb Miller, and David J. Lipman (1997),  
 "Gapped BLAST and PSI-BLAST: a new generation of protein database search programs", Nucleic Acids Res. 25:3389-3402.

Query: unkwown  
 (147 letters)

Database: prodom2010.1 multiple alignments  
 45,292,438 sequences; 2,147,483,647 total letters

ProDom domains producing High-scoring Segment Pairs:

Position	ProDom domain
Score E value	
19-147	#PD000577
678 3e-89	
42-134	#PDC1J6J2
180 2e-14	
45-145	#PDC5E2N9
202 1e-17	

>**PD000577** (Closest domain: LYSC\_CHICK 19-147)  
 Number of domains in family: 566  
 Commentary (automatic):  
 DISULFIDE BOND SUBNAME: FULL=LYSOZYME C RECNAME: ALTNAME:

EC=3.2.1.17 GLYCOSIDASE PRECURSOR  
Length = 129  
Score = 678 (265.8 bits), Expect = 3e-89  
Identities = 129/129 (100%), Positives = 129/129 (100%)

Query: 19  
KVFGRCELAAAMKRHGLDNYRGYSLGNWVCAAKFESNFNTQATNRNTDGSTDYGILO  
INS 78

KVFGRCELAAAMKRHGLDNYRGYSLGNWVCAAKFESNFNTQATNRNTDGSTDYGILO  
INS

Sbjct: 19  
KVFGRCELAAAMKRHGLDNYRGYSLGNWVCAAKFESNFNTQATNRNTDGSTDYGILO  
INS 78

Query: 79  
RWWCNDGRTPGSRNLCNIPCSALLSSDITASVNCAKKIVSDGNGMNAWVAWRNRCKG  
TDV 138

RWWCNDGRTPGSRNLCNIPCSALLSSDITASVNCAKKIVSDGNGMNAWVAWRNRCKG  
TDV

Sbjct: 79  
RWWCNDGRTPGSRNLCNIPCSALLSSDITASVNCAKKIVSDGNGMNAWVAWRNRCKG  
TDV 138

Query: 139 QAWIRGCRL 147  
QAWIRGCRL  
Sbjct: 139 QAWIRGCRL 147

>**PDC5E2N9** (Closest domain: Q7Q6R2\_ANOGA 366-501)  
Number of domains in family: 2  
Commentary (automatic):  
SUBNAME: DISULFIDE BOND FULL=LYSOZYME FULL=AGAP005717-PA  
REFERENCE C-6  
Length = 136  
Score = 202 (82.4 bits), Expect = 1e-17  
Identities = 44/107 (41%), Positives = 59/107 (55%), Gaps  
= 9/107 (8%)

Query: 45 NWVCAAKFESNFNTQATNR-  
NTDGSTDYGILQINSRWWCNDGRTPGSRNLCNIPCSALLS 103  
WVC A ES FNT A R N DGS D+G+ QI+ +WC+  
PG+ C + C AL  
Sbjct: 366 TWVCIAYHESRFNTSAEGRNLNADGSGDHGLFQISDIYWCS--  
-PPGNGWACGVSCDALKD 422

Query: 104 SDITASVNCAKKIVSD-----

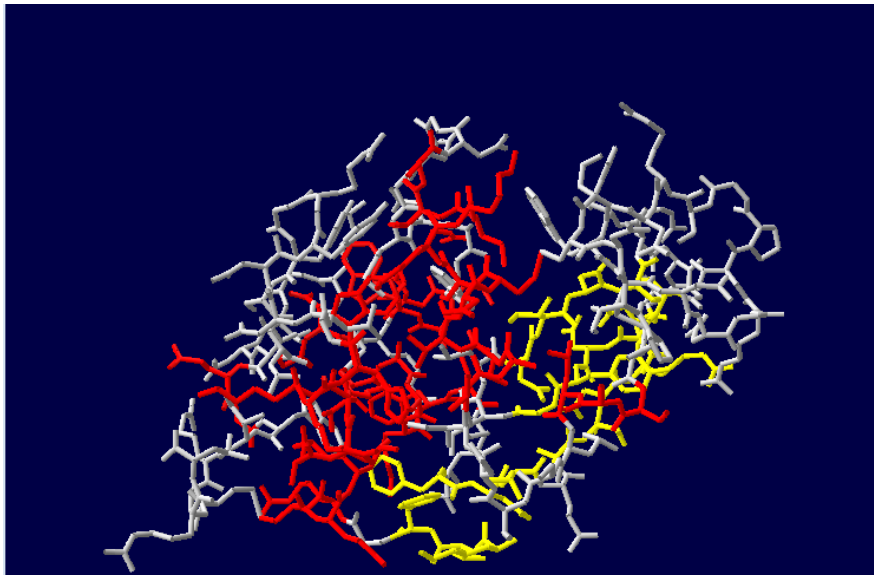
GNGMNAWVAWRNRCKGTDVQAWIRGC 145  
SDI+ V C K I + G+G NAW ++ C+ V  
++RGC  
Sbjct: 423  
SDISDDVQCVKTIYEEHQRLSGDGFNAWSVYKPYCQRDAVDTFVRGC 469

>**PDC1J6J2** (Closest domain: B4IWP9\_DROGR 949-1127)  
Number of domains in family: 3  
Commentary (automatic):  
SUBNAME: DISULFIDE BOND FULL=GI13315 FULL=GH16162  
Length = 179  
Score = 180 (73.9 bits), Expect = 2e-14  
Identities = 38/100 (38%), Positives = 55/100 (55%), Gaps  
= 10/100 (10%)

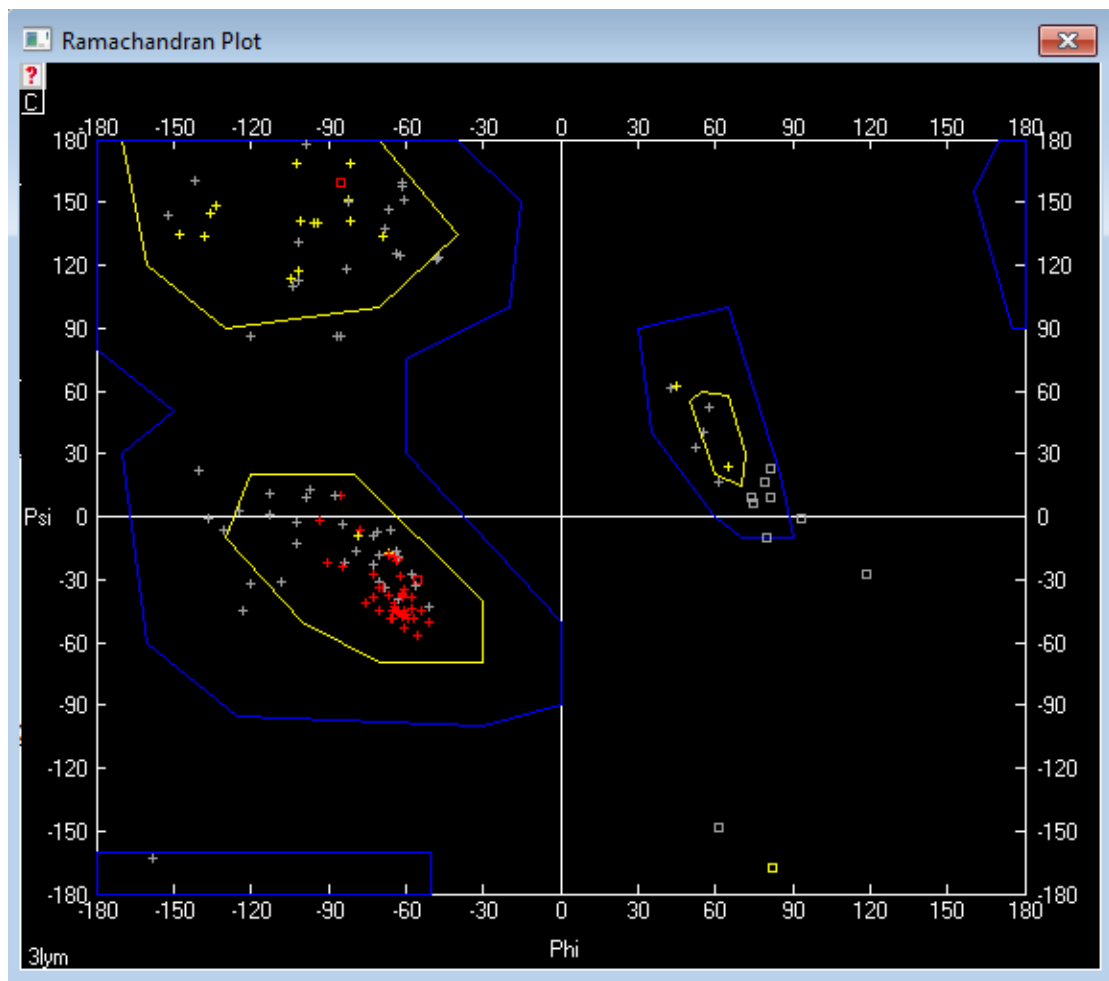
Query: 42 SLGNWVCAAKFESNFNTQATNR-  
NTDGSTDYGILQINSRWWC-NDGRTPGSRNLCNIPCS 99  
+ WVC A+ ES+++T A R NTD S D+G+ QI+ +WC  
+DG S C+I C  
Sbjct: 960  
EIPTWVCIAQHESYSTAAVGRNLNTDSSSEDHGLFQISDLYWCTHDG---  
SSGKACHIECD 1016

Query: 100 ALLSSDITASVNC AKKIVSD-----GNGMNAWVAWRNRCK  
134  
LL SDI+ + C K I + G+G AW + C+  
Sbjct: 1017 RLLDSDISDDIECIKTIYKEHTRISGDGFTAWTVYNGHCR  
1056

- Rendered by 3D/colored by secondary structure



- Ramachandran Plot



Τα αμινοξέα που εμφανίζονται εκτός των επιτρεπτών στερεοτακτικών δομών είναι τα:

**GLY102 GLY104 GLY 54**

Επειδή η πλευρική αλυσίδα της γλυκίνης είναι μικρή (μόλις ένα άτομο υδρογόνου), η γλυκίνη μπορεί να συμμετέχει σε πολλές διαμορφώσεις μη αποδεκτές για άλλα αμινοξέα