



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 01:43 PM UTC

PDB ID : 2EQL / pdb_00002eql
Title : CRYSTALLOGRAPHIC STUDIES OF A CALCIUM BINDING LYSOZYME
FROM EQUINE MILK AT 2.5 ANGSTROMS RESOLUTION
Authors : Tsuge, H.; Ago, H.; Miyano, M.
Deposited on : 1994-05-27
Resolution : 2.50 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4-5-2 with Phenix2.0
Xtrriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

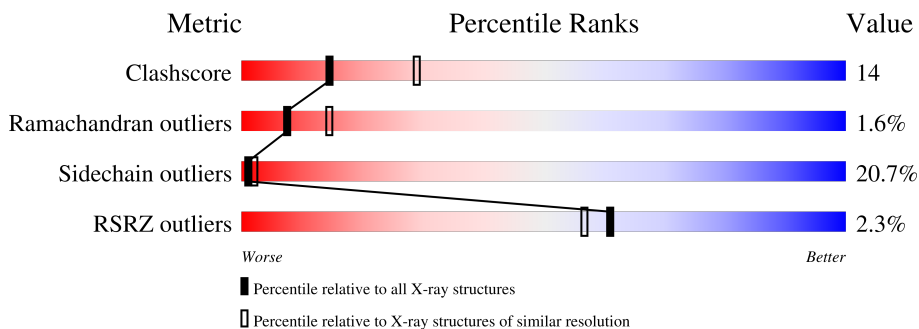
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	129	

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 1023 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

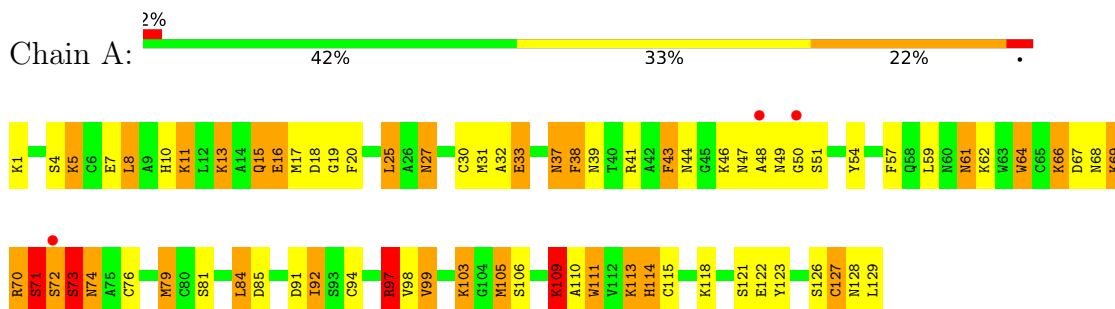
- Molecule 1 is a protein called HORSE MILK LYSOZYME.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace	
			Total	C	N	O				S
1	A	129	1023	636	180	195	12	0	0	0

3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

- Molecule 1: HORSE MILK LYSOZYME



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	54.10Å 57.20Å 38.70Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	7.00 – 2.50 7.00 – 2.50	Depositor EDS
% Data completeness (in resolution range)	(Not available) (7.00-2.50) 92.8 (7.00-2.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$	-	Xtrriage
Refinement program	X-PLOR	Depositor
R, R_{free}	0.234 , (Not available) 0.227 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	17.3	Xtrriage
Anisotropy	0.108	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 78.7	EDS
L-test for twinning ¹	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	1023	wwPDB-VP
Average B, all atoms (Å ²)	12.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 9.04% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality i

5.1 Standard geometry i

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	1.31	3/1044 (0.3%)	2.55	72/1399 (5.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	114	HIS	CD2-NE2	-5.97	1.31	1.37
1	A	97	ARG	CZ-NH1	5.45	1.40	1.32
1	A	8	LEU	CB-CG	-5.02	1.43	1.53

All (72) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	128	ASN	CA-CB-CG	-14.73	97.87	112.60
1	A	47	ASN	OD1-CG-ND2	-12.94	109.67	122.60
1	A	67	ASP	CA-CB-CG	12.11	124.71	112.60
1	A	127	CYS	CA-C-N	-11.26	107.71	122.79
1	A	127	CYS	C-N-CA	-11.26	107.71	122.79
1	A	16	GLU	CA-CB-CG	9.14	132.37	114.10
1	A	43	PHE	CA-CB-CG	-8.83	104.97	113.80
1	A	48	ALA	CA-C-O	8.24	129.15	120.42
1	A	97	ARG	NH1-CZ-NH2	-7.92	109.00	119.30
1	A	74	ASN	OD1-CG-ND2	-7.89	114.71	122.60
1	A	18	ASP	CA-C-O	-7.75	112.24	120.92
1	A	49	ASN	O-C-N	-7.63	115.26	123.42
1	A	33	GLU	CA-CB-CG	7.54	129.18	114.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	4	SER	N-CA-C	-7.47	100.89	110.53
1	A	85	ASP	N-CA-C	7.46	121.43	110.17
1	A	71	SER	CA-C-N	7.43	134.60	121.75
1	A	71	SER	C-N-CA	7.43	134.60	121.75
1	A	68	ASN	CB-CG-ND2	7.42	127.53	116.40
1	A	27	ASN	CA-CB-CG	-7.36	105.24	112.60
1	A	99	VAL	CB-CA-C	-7.34	99.25	111.29
1	A	73	SER	CA-C-O	-7.32	113.46	120.56
1	A	97	ARG	CA-CB-CG	7.13	128.37	114.10
1	A	38	PHE	CA-CB-CG	7.08	120.88	113.80
1	A	16	GLU	N-CA-C	6.84	121.00	112.24
1	A	15	GLN	CA-CB-CG	-6.64	100.82	114.10
1	A	54	TYR	N-CA-C	6.47	119.49	109.07
1	A	68	ASN	OD1-CG-ND2	-6.46	116.14	122.60
1	A	84	LEU	CD1-CG-CD2	-6.35	96.84	110.80
1	A	111	TRP	N-CA-C	-6.32	104.31	111.14
1	A	37	ASN	CA-CB-CG	-6.30	106.30	112.60
1	A	20	PHE	N-CA-C	6.29	119.11	110.24
1	A	127	CYS	CB-CA-C	-6.26	102.55	111.76
1	A	73	SER	N-CA-C	-6.21	99.66	108.23
1	A	5	LYS	CA-C-O	6.18	127.31	120.82
1	A	48	ALA	CA-C-N	6.18	132.93	121.62
1	A	48	ALA	C-N-CA	6.18	132.93	121.62
1	A	13	LYS	O-C-N	6.00	128.48	122.12
1	A	85	ASP	N-CA-CB	-5.90	101.01	110.51
1	A	61	ASN	CA-CB-CG	5.88	118.48	112.60
1	A	15	GLN	N-CA-C	5.83	120.10	112.92
1	A	69	LYS	N-CA-C	5.82	119.12	109.46
1	A	44	ASN	CA-CB-CG	-5.78	106.82	112.60
1	A	91	ASP	CA-CB-CG	5.78	118.38	112.60
1	A	15	GLN	N-CA-CB	-5.76	102.14	110.56
1	A	7	GLU	CA-CB-CG	5.74	125.57	114.10
1	A	11	LYS	N-CA-C	5.71	118.28	111.71
1	A	8	LEU	N-CA-CB	-5.70	101.73	110.16
1	A	79	MET	N-CA-C	-5.70	100.70	109.76
1	A	128	ASN	O-C-N	5.68	129.70	122.65
1	A	49	ASN	N-CA-C	-5.66	100.13	108.79
1	A	64	TRP	CG-CD1-NE1	-5.65	102.85	110.20
1	A	15	GLN	CA-C-N	5.60	130.53	122.40
1	A	15	GLN	C-N-CA	5.60	130.53	122.40
1	A	61	ASN	CA-C-O	-5.59	113.55	119.97
1	A	39	ASN	CA-CB-CG	5.57	118.17	112.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	69	LYS	CA-CB-CG	5.56	125.22	114.10
1	A	70	ARG	CA-CB-CG	5.53	125.17	114.10
1	A	126	SER	N-CA-CB	-5.53	102.45	110.47
1	A	32	ALA	CA-C-O	-5.48	114.71	120.63
1	A	64	TRP	CE2-CD2-CG	-5.42	100.70	107.20
1	A	74	ASN	N-CA-C	5.31	119.28	109.24
1	A	16	GLU	CB-CG-CD	5.27	121.56	112.60
1	A	109	LYS	N-CA-C	5.24	117.67	111.33
1	A	5	LYS	N-CA-CB	-5.21	102.45	110.01
1	A	114	HIS	CB-CG-CD2	-5.21	124.43	131.20
1	A	105	MET	CA-CB-CG	-5.19	103.71	114.10
1	A	17	MET	CG-SD-CE	-5.19	89.48	100.90
1	A	111	TRP	CG-CD1-NE1	-5.18	103.47	110.20
1	A	19	GLY	CA-C-N	5.16	128.17	120.90
1	A	19	GLY	C-N-CA	5.16	128.17	120.90
1	A	7	GLU	CB-CA-C	-5.12	102.61	110.81
1	A	64	TRP	N-CA-C	5.09	119.76	113.50

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	97	ARG	Sidechain

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1023	0	975	28	0
All	All	1023	0	975	28	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 14.

All (28) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:81:SER:HA	1:A:84:LEU:HD12	1.50	0.92
1:A:27:ASN:HB3	1:A:105:MET:HE1	1.69	0.75
1:A:43:PHE:HZ	1:A:84:LEU:HA	1.55	0.72
1:A:62:LYS:O	1:A:73:SER:HB2	1.93	0.67
1:A:33:GLU:HB3	1:A:38:PHE:CE1	2.31	0.65
1:A:127:CYS:HA	1:A:129:LEU:HD13	1.81	0.63
1:A:94:CYS:O	1:A:98:VAL:HG23	2.00	0.61
1:A:5:LYS:HD3	1:A:127:CYS:SG	2.42	0.60
1:A:66:LYS:HE3	1:A:79:MET:SD	2.45	0.57
1:A:64:TRP:O	1:A:76:CYS:HB2	2.05	0.56
1:A:109:LYS:HZ3	1:A:113:LYS:HD3	1.72	0.54
1:A:33:GLU:HB3	1:A:38:PHE:CZ	2.42	0.54
1:A:62:LYS:HG2	1:A:71:SER:HB3	1.92	0.52
1:A:61:ASN:O	1:A:72:SER:HA	2.11	0.51
1:A:110:ALA:O	1:A:114:HIS:HB2	2.12	0.49
1:A:43:PHE:CE2	1:A:84:LEU:HD23	2.47	0.49
1:A:27:ASN:HB3	1:A:105:MET:CE	2.41	0.49
1:A:81:SER:CA	1:A:84:LEU:HD12	2.34	0.48
1:A:66:LYS:HB2	1:A:74:ASN:ND2	2.30	0.47
1:A:31:MET:HE1	1:A:57:PHE:CD1	2.51	0.46
1:A:43:PHE:CZ	1:A:84:LEU:HD23	2.52	0.45
1:A:30:CYS:HA	1:A:123:TYR:OH	2.18	0.44
1:A:13:LYS:HB2	1:A:25:LEU:HD11	1.98	0.44
1:A:111:TRP:CD1	1:A:115:CYS:HB2	2.52	0.44
1:A:97:ARG:HE	1:A:97:ARG:HB2	1.71	0.43
1:A:15:GLN:OE1	1:A:92:ILE:HG21	2.19	0.42
1:A:15:GLN:O	1:A:16:GLU:HG3	2.20	0.42
1:A:103:LYS:O	1:A:106:SER:HB2	2.21	0.41

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	127/129 (98%)	115 (91%)	10 (8%)	2 (2%)	7 14

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	70	ARG
1	A	50	GLY

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	A	111/111 (100%)	88 (79%)	23 (21%)	1 2

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	LYS
1	A	8	LEU
1	A	10	HIS
1	A	11	LYS
1	A	25	LEU
1	A	37	ASN
1	A	41	ARG
1	A	46	LYS
1	A	51	SER
1	A	59	LEU
1	A	66	LYS
1	A	69	LYS
1	A	71	SER
1	A	72	SER
1	A	73	SER
1	A	92	ILE
1	A	99	VAL
1	A	103	LYS
1	A	109	LYS

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Mol	Chain	Res	Type
1	A	113	LYS
1	A	118	LYS
1	A	121	SER
1	A	122	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	47	ASN
1	A	58	GLN
1	A	77	ASN
1	A	114	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled '#RSRZ > 2' contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled 'Q < 0.9' lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	129/129 (100%)	-0.25	3 (2%) 61 57	2, 11, 24, 33	0

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	48	ALA	3.3
1	A	50	GLY	2.3
1	A	72	SER	2.1

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.