

Full wwPDB X-ray Structure Validation Report (i)

May 23, 2020 – 11:58 pm BST

PDB ID : 139L

> Title RAPID CRYSTALLIZATION OF T4 LYSOZYME BY INTERMOLECULAR

> > DISULFIDE CROSSLINKING

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Deposited on : 1993-09-01

1.70 Å(reported) Resolution

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 (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

MolProbity 4.02b-467

> 1.8.5 (274361), CSD as541be (2020) Mogul

Xtriage (Phenix) 1.13 EDS 2.11

Percentile statistics 20191225.v01 (using entries in the PDB archive December 25th 2019)

> Refmac 5.8.0158

7.0.044 (Gargrove) CCP4 Engh & Huber (2001)

Ideal geometry (proteins) Ideal geometry (DNA, RNA) Parkinson et al. (1996)

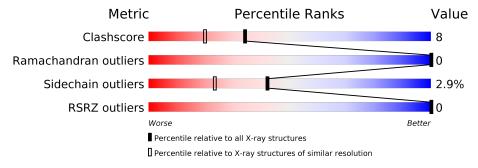
Validation Pipeline (wwPDB-VP) 2.11

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 1.70 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	Similar resolution	
Metric	$(\# \mathrm{Entries})$	$(\# \text{Entries, resolution range}(\text{\AA}))$	
Clashscore	141614	4695 (1.70-1.70)	
Ramachandran outliers	138981	4610 (1.70-1.70)	
Sidechain outliers	138945	4610 (1.70-1.70)	
RSRZ outliers	127900	4222 (1.70-1.70)	

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 on 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 <=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	164	80%	15%	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	CL	A	178	_	-	_	X



2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 1447 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 T4 LYSOZYME.

Mol	Chain	Residues		At	oms			ZeroOcc	AltConf	Trace
1	A	162	Total 1291	C 813	N 234	O 237	S 7	0	0	0

There are 4 discrepancies between the modelled and reference sequences:

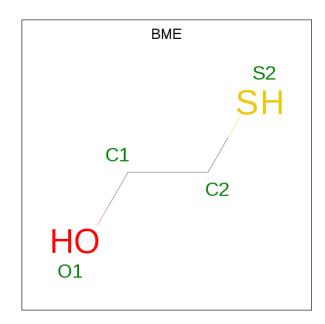
Chain	Residue	Modelled	Actual	Comment	Reference
A	54	THR	CYS	CONFLICT	UNP P00720
A	68	CYS	ASN	CONFLICT	UNP P00720
A	93	CYS	ALA	CONFLICT	UNP P00720
A	97	ALA	CYS	CONFLICT	UNP P00720

• Molecule 2 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	2	Total Cl 2 2	0	0

• Molecule 3 is BETA-MERCAPTOETHANOL (three-letter code: BME) (formula: C₂H₆OS).





$\overline{\text{Mol}}$	Chain	Residues	${f Atoms}$				ZeroOcc	AltConf
3	A	1	Total 4	C 2	O 1	S 1	0	0
3	A	1	Total 4	C 2	O 1	S 1	0	0

• Molecule 4 is water.

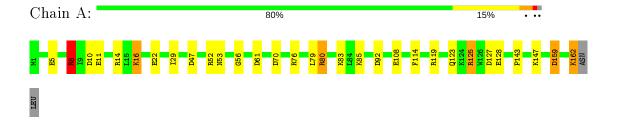
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	146	Total O 146 146	0	0



3 Residue-property plots (i)

These plots are drawn for all protein, RNA and DNA 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: T4 LYSOZYME





4 Data and refinement statistics (i)

Property	Value	Source			
Space group	P 32 2 1	Depositor			
Cell constants	$60.86 \text{\AA} 60.86 \text{Å} 97.42 \text{Å}$	Donositor			
a, b, c, α , β , γ	90.00° 90.00° 120.00°	Depositor			
Resolution (Å)	(Not available) – 1.70	Depositor			
Resolution (A)	25.81 - 1.70	Depositor Depositor EDS Depositor EDS Depositor Depositor Xtriage Depositor Depositor Very Striage Television Striage Televisio			
% Data completeness	(Not available) ((Not available)-1.70)	Depositor			
(in resolution range)	82.9 (25.81-1.70)	EDS			
R_{merge}	(Not available)	Depositor			
R_{sym}	(Not available)	Depositor			
$< I/\sigma(I) > 1$	1.24 (at 1.70Å)	Xtriage			
Refinement program	TNT	Depositor			
R, R_{free}	0.164 , (Not available)	Depositor			
It, It free	0.163 , (Not available)	DCC			
R_{free} test set	No test flags present.	wwPDB-VP			
Wilson B-factor (Å ²)	14.1	Xtriage			
Anisotropy	0.128	Xtriage			
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.34, 92.1	EDS			
L-test for twinning ²	$< L >=0.48, < L^2>=0.31$	Xtriage			
Estimated twinning fraction	0.038 for -h,-k,l	Xtriage			
F_o, F_c correlation	0.97	EDS			
Total number of atoms	1447	wwPDB-VP			
Average B, all atoms (Å ²)	23.0	wwPDB-VP			

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 8.31% of the height of the origin peak. No significant pseudotranslation is detected.

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



¹Intensities estimated from amplitudes.

5 Model quality (i)

5.1 Standard geometry (i)

Bond lengths and bond angles in the following residue types are not validated in this section: CL, BME

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	Boı	nd lengths	В	ond angles
WIOI	Chain	RMSZ	# Z > 5	RMSZ	# Z >5
1	A	2.94	9/1311 (0.7%)	1.57	20/1765~(1.1%)

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$\operatorname{Observed}(\operatorname{\AA})$	$Ideal(\AA)$
1	A	162	LYS	N-CA	64.04	2.74	1.46
1	A	162	LYS	CA-C	59.53	3.07	1.52
1	A	162	LYS	C-O	47.93	2.14	1.23
1	A	11	GLU	CD-OE1	6.58	1.32	1.25
1	A	128	GLU	CD-OE1	6.43	1.32	1.25
1	A	11	GLU	CD-OE2	-6.22	1.18	1.25
1	A	108	GLU	CD-OE1	6.01	1.32	1.25
1	A	22	GLU	CD-OE1	5.69	1.31	1.25
1	A	5	GLU	CD-OE2	5.47	1.31	1.25

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^o)$	$\operatorname{Ideal}({}^o)$
1	A	162	LYS	CA-C-O	-28.08	61.13	120.10
1	A	162	LYS	N-CA-C	-20.05	56.87	111.00
1	A	47	ASP	CB-CG-OD2	-11.91	107.58	118.30
1	A	127	ASP	CB-CG-OD1	10.04	127.33	118.30
1	A	47	ASP	CB-CG-OD1	8.93	126.34	118.30
1	A	80	ARG	NE-CZ-NH1	8.66	124.63	120.30
1	A	127	ASP	CB-CG-OD2	-8.01	111.09	118.30
1	A	70	ASP	CB-CG-OD1	7.24	124.81	118.30
1	A	8	ARG	NE-CZ-NH1	6.61	123.61	120.30
1	A	92	ASP	CB-CG-OD2	6.48	124.14	118.30
1	A	70	ASP	CB-CG-OD2	-6.15	112.76	118.30
1	A	162	LYS	N-CA-CB	-6.08	99.66	110.60

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Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$\mathbf{Observed}(^o)$	$Ideal(^{o})$
1	A	80	ARG	NE-CZ-NH2	-5.95	117.33	120.30
1	A	92	ASP	CB-CG-OD1	-5.80	113.08	118.30
1	A	125	ARG	NE-CZ-NH1	5.63	123.12	120.30
1	A	10	ASP	CB-CG-OD1	5.63	123.36	118.30
1	A	10	ASP	CB-CG-OD2	-5.46	113.39	118.30
1	A	61	ASP	CB-CG-OD1	-5.32	113.52	118.30
1	A	159	ASP	CB-CG-OD1	-5.27	113.55	118.30
1	A	159	ASP	CB-CG-OD2	5.21	122.99	118.30

There are no chirality outliers.

There are no planarity outliers.

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	1291	0	1318	22	0
2	A	2	0	0	0	0
3	A	8	0	10	0	0
4	A	146	0	0	4	0
All	All	1447	0	1328	22	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 8.

All (22) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	$egin{array}{ll} ext{Interatomic} \ ext{distance} \ (ext{\AA}) \end{array}$	Clash overlap (Å)
1:A:162:LYS:O	1:A:162:LYS:C	2.14	0.85
1:A:8:ARG:HH11	1:A:8:ARG:HG3	1.46	0.80
1:A:114:PHE:HA	4:A:326:HOH:O	1.90	0.70
1:A:80:ARG:NH2	4:A:242:HOH:O	2.30	0.65
1:A:76:ARG:NH1	1:A:76:ARG:HB2	2.14	0.63
1:A:125:ARG:NE	4:A:317:HOH:O	2.31	0.58
1:A:76:ARG:O	1:A:80:ARG:HG3	2.04	0.57
1:A:8:ARG:NH1	1:A:8:ARG:HG3	2.20	0.54

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Atom-1	Atom-2	Interatomic	Clash
Atom-1	Atom-2	${f distance} \; ({f \AA})$	overlap (Å)
1:A:76:ARG:CB	1:A:76:ARG:HH11	2.20	0.53
1:A:143:PRO:O	1:A:147:LYS:HG3	2.11	0.51
1:A:162:LYS:N	1:A:162:LYS:CA	2.74	0.51
1:A:76:ARG:HB2	1:A:76:ARG:HH11	1.76	0.51
1:A:29:ILE:O	1:A:29:ILE:HG22	2.13	0.48
1:A:159:ASP:OD1	1:A:162:LYS:NZ	2.36	0.48
1:A:119:ARG:HG2	1:A:123:GLN:NE2	2.29	0.47
1:A:16:LYS:HE3	1:A:56:GLY:O	2.14	0.47
1:A:119:ARG:HG2	1:A:123:GLN:HE21	1.81	0.46
1:A:52:ARG:HG3	1:A:52:ARG:HH11	1.81	0.46
1:A:14:ARG:HG3	4:A:298:HOH:O	2.15	0.46
1:A:79:LEU:O	1:A:85:LYS:HD2	2.17	0.45
1:A:76:ARG:NH1	1:A:76:ARG:CB	2.80	0.43
1:A:8:ARG:O	1:A:8:ARG:HG3	2.17	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	160/164 (98%)	157 (98%)	3 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	136/138 (99%)	132 (97%)	4 (3%)	42 23

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	8	ARG
1	A	16	LYS
1	A	53	ASN
1	A	83	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	69	GLN
1	A	123	GLN
1	A	140	ASN

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 carbohydrates in this entry.

5.6 Ligand geometry (i)

Of 4 ligands modelled in this entry, 2 are monoatomic - leaving 2 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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| > 2 is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).



Mol	Т	Chain	Dog	T in le	B	ond leng	$_{ m gths}$	В	ond ang	gles
MIOI	Type	Chain	Res	Link	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	BME	A	902	3	3,3,3	0.82	0	1,2,2	0.25	0
3	BME	A	901	3	3,3,3	0.10	0	1,2,2	1.04	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	${f Res}$	Link	Chirals	Torsions	Rings
3	BME	A	902	3	-	1/1/1/1	-
3	BME	A	901	3	-	1/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	902	BME	O1-C1-C2-S2
3	A	901	BME	O1-C1-C2-S2

There are no ring outliers.

No monomer is involved in short contacts.

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, 95^{th} 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 $>$	$\#\text{RSRZ}{>}2$	$OWAB(A^2)$	Q < 0.9
1	A	162/164 (98%)	-0.73	0 100 100	7, 16, 48, 70	0

There are no RSRZ outliers to report.

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 carbohydrates in this entry.

6.4 Ligands (i)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	${f Res}$	Atoms	RSCC	RSR	${f B-factors(\AA^2)}$	Q<0.9
2	CL	A	178	1/1	0.33	0.68	32,32,32,32	1
3	BME	A	902	4/4	0.98	0.08	16,23,24,29	0
3	BME	A	901	4/4	0.98	0.12	26,27,29,39	0
2	CL	A	173	1/1	0.99	0.27	1,1,1,1	0

6.5 Other polymers (i)

There are no such residues in this entry.

