

# Full wwPDB X-ray Structure Validation Report (i)

May 25, 2020 – 06:02 pm BST

PDB ID : 1LJ3

Title : CRYSTAL STRUCTURE OF MONOCLINIC LYSOZYME GROWN AT PH

4.6

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Deposited on : 2002-04-19

Resolution : 2.00 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 (i) symbol.

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

MolProbity : 4.02b-467

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

Xtriage (Phenix) : 1.13 EDS : 2.11

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

Refmac : 5.8.0158

CCP4 : 7.0.044 (Gargrove) roteins) : Engh & Huber (2001)

Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)

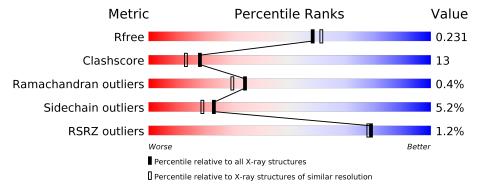
Validation Pipeline (wwPDB-VP) : 2.11

## 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 2.00 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	$\begin{array}{c} \text{Whole archive} \\ (\#\text{Entries}) \end{array}$	$\begin{array}{c} {\rm Similar\ resolution} \\ (\#{\rm Entries,\ resolution\ range(\AA)}) \end{array}$
$R_{free}$	130704	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (2.00-2.00)

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	129	76%	21%	•
1	В	129	78%	19%	•



## 2 Entry composition (i)

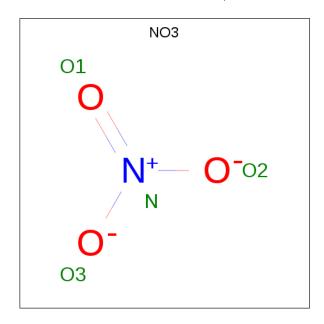
There are 3 unique types of molecules in this entry. The entry contains 2301 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Lysozyme C.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
1	А	129	Total	0	N	О	S	0	0	0
	71	120	1001	613	193	185	10		O	0
1	D	129	Total	С	N	Ο	S	0	0	0
1	Б	129	1001	613	193	185	10	0	U	0

• Molecule 2 is NITRATE ION (three-letter code: NO3) (formula: NO<sub>3</sub>).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total N O 4 1 3	0	0
2	A	1	Total N O 4 1 3	0	0
2	В	1	Total N O 4 1 3	0	0
2	В	1	Total N O 4 1 3	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf	
2	В	1	Total 4	N 1	O 3	0	0

### • Molecule 3 is water.

Mol	Chain	Residues	${f Atoms}$	ZeroOcc	AltConf
3	A	147	Total O 147 147	0	0
3	В	132	Total O 132 132	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.



## 4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants	$27.89 \text{\AA} - 62.76 \text{Å} - 60.27 \text{Å}$	D:4
a, b, c, $\alpha$ , $\beta$ , $\gamma$	$90.00^{\circ}$ $90.91^{\circ}$ $90.00^{\circ}$	Depositor
Resolution (Å)	30.00 - 2.00	Depositor
Resolution (A)	27.88 - 2.00	EDS
% Data completeness	94.4 (30.00-2.00)	Depositor
(in resolution range)	94.6 (27.88-2.00)	EDS
$R_{merge}$	(Not available)	Depositor
$R_{sum}$	(Not available)	Depositor
$< I/\sigma(I) > 1$	5.80 (at 1.99Å)	Xtriage
Refinement program	CNS 0.4	Depositor
D D.	0.178 , 0.233	Depositor
$R, R_{free}$	0.176 , $0.231$	DCC
$R_{free}$ test set	1078 reflections (8.08%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	20.7	Xtriage
Anisotropy	0.379	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$ , $B_{sol}(Å^2)$	0.32, 46.0	EDS
L-test for twinning <sup>2</sup>	$< L >=0.50, < L^2>=0.34$	Xtriage
	0.014 for -h,-l,-k	
Estimated twinning fraction	0.000 for $-h,l,k$	Xtriage
	0.048  for  h,-k,-l	
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	2301	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The analyses of the Patterson function reveals a significant off-origin peak that is 31.09 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.1720e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.

<sup>&</sup>lt;sup>2</sup>Theoretical 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.



<sup>&</sup>lt;sup>1</sup>Intensities 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: NO3

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	Bo	nd angles
MIOI	Chain	RMSZ	# Z >5	RMSZ	# Z  > 5
1	A	0.31	0/1021	0.60	1/1379 (0.1%)
1	В	0.30	0/1021	0.64	0/1379
All	All	0.30	0/2042	0.62	1/2758 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	${f Z}$	$\mathbf{Observed}(^o)$	$\operatorname{Ideal}({}^o)$
1	A	4	GLY	N-CA-C	-5.17	100.18	113.10

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	1001	0	967	23	0
1	В	1001	0	959	26	0
2	A	8	0	0	0	0
2	В	12	0	0	0	0
3	A	147	0	0	4	0
3	В	132	0	0	3	0
All	All	2301	0	1926	49	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 13.

All (49) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic	Clash
Atom-1	Atom-2	${ m distance} \; ({ m \AA})$	overlap (Å)
1:B:77:ASN:HB2	3:B:492:HOH:O	1.77	0.84
1:B:115:CYS:O	1:B:118:THR:HG22	1.78	0.82
1:B:45:ARG:NH2	1:B:49:GLY:HA2	1.99	0.78
1:B:51:THR:HG21	1:B:68:ARG:HH21	1.56	0.70
1:B:45:ARG:HG2	1:B:45:ARG:HH11	1.64	0.63
1:A:14:ARG:HD2	3:A:493:HOH:O	2.01	0.60
1:A:45:ARG:HG2	3:A:284:HOH:O	2.02	0.59
1:A:124:ILE:HG23	1:A:124:ILE:O	2.05	0.57
1:B:45:ARG:HH22	1:B:49:GLY:HA2	1.69	0.57
1:B:63:TRP:O	1:B:76:CYS:HB2	2.04	0.55
1:B:97:LYS:HB3	1:B:97:LYS:NZ	2.21	0.55
1:B:8:LEU:HD22	1:B:12:MET:HG3	1.88	0.54
1:A:33:LYS:HD3	1:A:33:LYS:C	2.28	0.54
1:B:45:ARG:C	1:B:45:ARG:HD3	2.28	0.53
1:A:64:CYS:HG	1:A:80:CYS:HG	0.55	0.53
1:B:45:ARG:NH1	1:B:45:ARG:HG2	2.23	0.53
1:B:118:THR:HG21	3:B:389:HOH:O	2.10	0.52
1:A:33:LYS:HE3	1:A:37:ASN:OD1	2.10	0.51
1:A:61:ARG:HD3	1:A:69:THR:CG2	2.42	0.50
1:B:8:LEU:HD12	1:B:38:PHE:CD1	2.47	0.49
1:B:121:GLN:NE2	3:B:269:HOH:O	2.45	0.49
1:A:18:ASP:O	1:A:19:ASN:HB2	2.12	0.49
1:A:3:PHE:HD2	1:A:7:GLU:HG2	1.78	0.49
1:A:45:ARG:HG3	1:A:45:ARG:HH11	1.78	0.48
1:B:121:GLN:HA	1:B:121:GLN:HE21	1.78	0.47
1:A:58:ILE:HB	1:A:83:LEU:HD13	1.97	0.47
1:B:33:LYS:HG2	1:B:123:TRP:CH2	2.50	0.47
1:B:61:ARG:HD2	1:B:61:ARG:O	2.14	0.47
1:A:15:HIS:HB3	1:A:92:VAL:HG11	1.98	0.46
1:A:63:TRP:O	1:A:76:CYS:HB2	2.15	0.46
1:A:66:ASP:HB3	1:A:80:CYS:SG	2.55	0.46
1:A:124:ILE:CG2	1:A:124:ILE:O	2.63	0.46
1:B:55:ILE:HG23	1:B:56:LEU:HG	1.99	0.44
1:A:73:ARG:HD2	3:A:436:HOH:O	2.17	0.44
1:B:51:THR:HG21	1:B:68:ARG:NH2	2.28	0.44
1:B:62:TRP:CZ3	1:B:73:ARG:HD2	2.52	0.44
1:B:45:ARG:HD3	1:B:46:ASN:N	2.34	0.43

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Atom-1	Atom-2	Interatomic	Clash
Atom-1	Atom-2	$oxed{  ext{distance }(\mathring{\mathbf{A}}) }$	overlap (Å)
1:A:124:ILE:HG13	1:A:129:LEU:HD11	2.00	0.43
1:B:69:THR:O	1:B:72:SER:HB2	2.18	0.43
1:B:70:PRO:O	1:B:72:SER:N	2.51	0.43
1:A:88:ILE:HA	3:A:317:HOH:O	2.19	0.43
1:A:78:ILE:HD12	1:A:78:ILE:O	2.20	0.42
1:A:45:ARG:NH1	1:A:45:ARG:HG3	2.35	0.42
1:B:70:PRO:O	1:B:71:GLY:C	2.57	0.42
1:A:103:ASN:O	1:A:106:ASN:HB2	2.20	0.42
1:A:125:ARG:O	1:A:125:ARG:HG3	2.19	0.41
1:B:61:ARG:HA	1:B:69:THR:HG21	2.03	0.41
1:A:114:ARG:HG2	1:A:114:ARG:NH1	2.36	0.40
1:B:62:TRP:HZ3	1:B:73:ARG:HD2	1.87	0.40

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	Perce	entiles
1	A	127/129 (98%)	126 (99%)	1 (1%)	0	100	100
1	В	127/129 (98%)	123 (97%)	3 (2%)	1 (1%)	19	13
All	All	254/258 (98%)	249 (98%)	4 (2%)	1 (0%)	34	30

#### All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	В	71	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	Percentiles		
1	A	105/105 (100%)	99 (94%)	6 (6%)	20	16
1	В	$105/105 \; (100\%)$	100 (95%)	5 (5%)	25	22
All	All	210/210 (100%)	199 (95%)	11 (5%)	23	19

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	33	LYS
1	A	45	ARG
1	A	61	ARG
1	A	112	ARG
1	A	124	ILE
1	A	128	ARG
1	В	8	LEU
1	В	45	ARG
1	В	51	THR
1	В	84	LEU
1	В	103	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	${f Res}$	$\mathbf{Type}$
1	A	19	ASN
1	A	41	GLN
1	В	41	GLN
1	В	103	ASN
1	В	113	ASN
1	В	121	GLN

#### 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)

5 ligands are modelled in this entry.

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 Type		Chain	Chain	Chain	Res	Link	Bond lengths			Bond angles		
WIOI	Type	nes		Lilik	Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2		
2	NO3	A	703	-	1,3,3	0.12	0	0,3,3	0.00	-		
2	NO3	В	700	-	1,3,3	0.24	0	0,3,3	0.00	-		
2	NO3	В	701	-	1,3,3	0.17	0	0,3,3	0.00	-		
2	NO3	В	704	-	1,3,3	0.18	0	0,3,3	0.00	-		
2	NO3	A	702	-	1,3,3	0.25	0	0,3,3	0.00	-		

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

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 $>$	# RSRZ > 2		$\mathbf{OWAB}(\mathbf{\mathring{A}}^2)$	Q<0.9
1	A	129/129 (100%)	-0.22	1 (0%) 86	85	13, 22, 34, 42	0
1	В	$129/129 \ (100\%)$	-0.25	2 (1%) 72	70	13, 20, 35, 39	0
All	All	258/258 (100%)	-0.24	3 (1%) 79	78	13, 21, 34, 42	0

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	73	ARG	4.2
1	В	72	SER	3.2
1	В	45	ARG	3.1

## 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 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	Res	Atoms	RSCC	RSR	${f B\text{-factors}}({f \AA}^2)$	Q < 0.9
2	NO3	A	703	4/4	0.85	0.27	73,73,73,73	0
2	NO3	В	704	4/4	0.92	0.19	52,53,53,53	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	$\operatorname{\textbf{B-factors}}(\AA^2)$	Q<0.9
2	NO3	В	701	4/4	0.94	0.11	39,39,39,40	0
2	NO3	A	702	4/4	0.94	0.14	39,39,39,39	0
2	NO3	В	700	4/4	0.96	0.09	27,28,28,28	0

## 6.5 Other polymers (i)

There are no such residues in this entry.

